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<b>(21) International Application Number:</b> PCT/US92/08788 <b>(22) International Filing Date:</b> 15 October 1992 (15.10.92) <b>(30) Priority data:</b> 781,477 18 October 1991 (18.10.91) US 866,931 10 April 1992 (10.04.92) US <b>(60) Parent Application or Grant</b> <b>(63) Related by Continuation</b> US 866,931 (CIP) Filed on 10 April 1992 (10.04.92) <b>(71) Applicant (for all designated States except US):</b> GENENTECH, INC. [US/US]; 460 Point San Bruno Boulevard, South San Francisco, CA 94080-4990 (US). <b>(72) Inventors; and</b> <b>(75) Inventors/Applicants (for US only):</b> BLACKBURN, Brent [US/US]; 11 Hillview Court, San Francisco, CA 94124 (US). BARKER, Peter [US/US]; 357 Avenue Del Oro, El Granada, CA 94018 (US). GADEK, Thomas [US/US]; 2838 Chelsea Drive, Oakland, CA 94611 (US). McDOWELL, Robert [US/US]; 11 Hillview Court, San Francisco, CA 94124 (US). McGEE, Lawrence [US/US]; 1 Crater Lake Way, Pacifica, CA 94044 (US). SOMERS, Todd [US/US]; P.O. Box 370315, Montara, CA 94037 (US). WEBB, Rob [US/US]; 802 Stetson Street, Moss Beach, CA 94035 (US). ROBARGE, Kirk [US/US]; 1260 Haight Street, #4, San Francisco, CA 94117 (US).		<b>(74) Agents:</b> WINTER, Daryl, B. et al.; Genentech, Inc., 460 Point San Bruno Boulevard, South San Francisco, CA 94080-4990 (US). <b>(81) Designated States:</b> AU, BB, BG, BR, CA, CS, DK, FI, HU, JP, KP, KR, LK, MG, MN, MW, NO, PL, RO, RU, SD, US, European patent (AT, BE, CH, DE, DK, ES, FR, GB, GR, IE, IT, LU, MC, NL, SE), OAPI patent (BF, BJ, CF, CG, CI, CM, GA, GN, ML, MR, SN, TD, TG). <b>Published</b> <i>With international search report.</i>
<b>(54) Title:</b> NONPEPTIDYL INTEGRIN INHIBITORS HAVING SPECIFICITY FOR THE GPII <sub>b</sub> III <sub>a</sub> RECEPTOR  <b>(57) Abstract</b>  A benzodiazepinedione derivative which acts as a nonpeptidyl platelet aggregation inhibitor is provided. This inhibitor potently inhibits fibrinogen binding to the GPII <sub>b</sub> III <sub>a</sub> receptor and is provided in therapeutic compositions for the treatment of diseases for which blocking platelet aggregation is indicated. These nonpeptidyl inhibitors are provided in combination with thrombolytics and anticoagulants.		

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## FIELD OF THE INVENTION

5           The present invention relates to nonpeptidyl inhibitors of integrin receptors. Specifically, the invention is directed to antagonists of the final common pathway of platelet aggregation. These antagonists act as potent antithrombotics. The invention further relates to therapeutic applications of these nonpeptidyl inhibitors in diseases for which blocking platelet aggregation is indicated.

## 10 BACKGROUND OF THE INVENTION

Platelets are particles found in whole blood that initiate and provide the structural basis for the hemostatic plug necessary to stop bleeding. Platelets depend on adhesive interactions with extracellular proteins and other cells for proper function (see Hawiger, J. *Atherosclerosis Reviews* 21:165-186 (1990) and Roth J. R. *Immunology Today* 13(2):100-105 (1992)). The external platelet plasma membrane surface is covered with a variety of membrane bound glycoproteins, many of which have recognition and adhesive functions. Perhaps the most abundant platelet membrane adhesive proteins belong to the integrin superfamily which include the glycoproteins; GPIIb/IIIa, GPIa/IIa, GPIc/IIa, GPIb/IX, and the fibronectin and vitronectin receptors (Hynes, R. O., *Cell*, 48: 549 (1987). Each integrin receptor is an  $\alpha\beta$  heterodimer displaying characteristic affinity and specificity toward various protein ligands found in serum and/or the extracellular matrix including; von Willebrand factor (vWF), collagen, entactin, tenascin, fibronectin (Fn), vitronectin (Vn), and laminin, as well as fibrinogen (Fg) and thrombospondin (see Kieffer *et al.*, *Ann. Rev. Cell Biol.* 6:329-357(1990) and Ruoslahti, *J. Clin. Invest.*, 87:1-5 (1991)). The most abundant integrin found on the surface of normal platelets is GPIIb/IIIa comprising about 50,000 molecules per platelet and representing about 2% of the total platelet protein. GPIIb/IIIa is a non-covalent, calcium ion dependent heterodimeric complex (Jennings, *et al.*, *J. Biol. Chem.* 257: 10458 (1982)) that is restricted in distribution to platelets and other cells of the megakaryocytic lineage (Kieffer *et al.*, *supra*). On activated platelets, GPIIb/IIIa promiscuously binds a number of protein ligands with varying affinities, including; fibrinogen, fibronectin, von Willebrand factor, vitronectin and thrombospondin (Plow *et al.*, *Biochemistry of Platelets*, Phillips and Shuman eds., p. 225-256, Orlando: Academic Press [1986]). Each of these protein ligands contain at least one tripeptide sequence Arg-Gly-Asp (RGD) which is commonly referred to as the "recognition sequence". It is believed the most important interactions mediating platelet aggregation involve GPIIb/IIIa binding with the trimeric fibrinogen and, to a lesser extent, with the filamentous von Willebrand factor (Kieffer *et al.*, *supra* and Albeda *et al.*, *The FASEB Journal*, 4:2868-2880 [1990]).

GPIIb/IIIa binding to its natural ligands can be inhibited to varying degrees by peptides and proteins containing the amino acid recognition sequences; Arg-Gly-Asp (Ruoslahti, *supra* 40 and EPO 368 486, assigned to Merck & Co.), Lys-Gly-Asp (KGD), and the fibrinogen  $\gamma$ -chain

carboxy-terminal dodecapeptide HHLGGAKQAGDV and analogues thereof (Timmons *et al.*, *Biochemistry*, 28:2919-2922 [1989]).

#### B. The Hyperthrombotic State

Many common human disorders are characteristically associated with a hyperthrombotic state leading to intravascular thrombi and emboli. These are a major cause of medical morbidity, leading to infarction, stroke and phlebitis, and of mortality from stroke and pulmonary and cardiac emboli. Patients with atherosclerosis are predisposed to arterial thromboembolic phenomena for a variety of reasons. Atherosclerotic plaques form niduses or platelet plugs and thrombi that lead to vascular narrowing and occlusion, resulting in myocardial and cerebral ischemic disease. This may happen spontaneously or following procedures such as angioplasty or endarterectomy. Thrombi that break off and are released into the circulation cause infarction of different organs, especially the brain, extremities, heart and kidneys.

In addition to being involved in arterial thrombosis, platelets may also play a role in venous thrombosis. A large percentage of such patients have no antecedent risk factors and develop venous thrombophlebitis and subsequent pulmonary emboli without a known cause. Other patients who form venous thrombi have underlying diseases known to predispose them to these syndromes. Some of these patients may have genetic or acquired deficiencies of factors that normally prevent hypercoagulability, such as antithrombin-3. Others have mechanical obstructions to venous flow, such as tumor masses, that lead to low flow states and thrombosis. Patients with malignancy have a high incidence of thrombotic phenomena for unclear reasons. Antithrombotic therapy in this situation with currently available agents is dangerous and often ineffective.

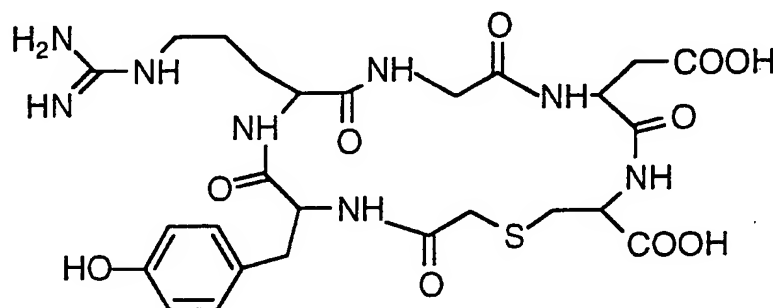
Patients whose blood flows over artificial surfaces, such as prosthetic synthetic cardiac valves or through extracorporeal perfusion devices, are also at risk for the development of platelet plugs, thrombi and emboli. It is standard practice that patients with artificial cardiac valves be treated chronically with anti-coagulants. However, in all instances, platelet activation and emboli formation may still occur despite adequate anticoagulation treatment.

Thus, a large category of patients, including those with atherosclerosis, coronary artery disease, artificial heart valves, cancer, and a history of stroke, phlebitis, or pulmonary emboli, are candidates for limited or chronic antithrombotic therapy. The number of available therapeutic agents is limited and these, for the most part, act by inhibiting or reducing levels of circulating clotting factors. These agents are frequently not effective against the patient's underlying hematologic problem, which often concerns an increased propensity for platelet aggregation and adhesion. They also cause the patient to be susceptible to abnormal bleeding. Available antiplatelet agents, such as aspirin, inhibit only part of the platelet activation process and are therefore often inadequate for therapy and also cause the patient to be susceptible to abnormal bleeding. An ideal anti-thrombotic drug would have many properties currently not available (see e.g. Sixma, *et al.* *Thrombosis Research* 67: 305-311 [1992]).

## C. Therapeutic Agents

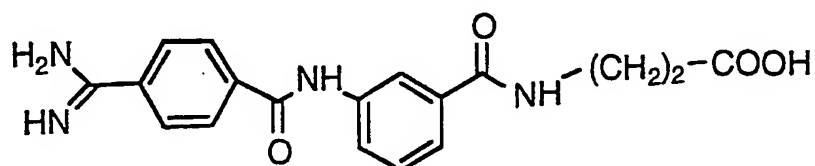
An agent which effectively inhibits the final common pathway of platelet activation, namely fibrinogen binding to the GP IIb/IIIa receptor, should accordingly be useful in a large group of disorders characterized by a hyperthrombotic state as described above.

- Such agents include anti-thrombotic peptides and pseudopeptides capable of inhibiting platelet aggregation. Ruoslahti *et al.* (U.S. patent no. 4,578,079) suggest that tetrapeptides containing the RGD sequence may be used to inhibit platelet aggregation. Zimmerman *et al.* (U.S. patent no. 4,683,291) disclose that positively charged amino acid residues (e. g. Arg and Lys) and homologues located before or toward the amino terminus of the RGD sequence are superior for inhibiting fibrinogen-platelet binding. Adams *et al.* (U.S. patent no. 4,857,508) describe superior results for *in vitro* inhibition of human platelet aggregation in platelet-rich plasma for linear tetrapeptides containing O-methyl-Tyr-amide immediately following the RGD (or homo-RGD) sequence. Barker *et al.*, WO 90/01331, disclose substantially rigid RGD cyclic peptides possessing high affinity for the GP IIb/IIIa receptor. A particularly efficacious rigid RGD cyclic peptide described by Barker *et al.* is represented by the following structure:



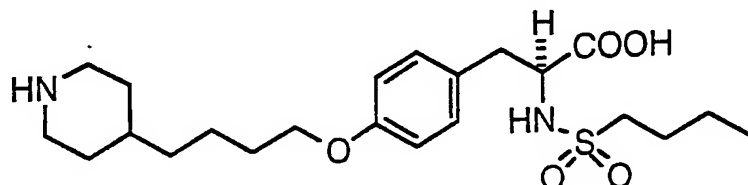
- Tjoeng *et al.* (U.S. patent no. 4,879,313) describe peptide mimetic platelet aggregation inhibitors in which the first two residues of the RGD sequence are replaced by the pseudodipeptidyl 8-guanidino-octanoyl moiety. Other peptidomimetics in which the Arg of the RGD sequence has been altered include; WO89/07609 (homo-Arg[Har]), EP 341 915 (Har and alkyl-Arg), WO90/15620 (Har and amidino derivatives e. g. imidazolyl, imidazolyl, and substituted imidazolyl), EP 422 937 (aryl-, arylalkyl-, and cycloalkyl-amines), WO91/07976 (alkylamidino and alkylamino derivatives), and WO91/04247 (alkylamino and alkylguanidino proline derivatives). See also EP 384 362 (glycine derivatives) and EP 381 033.

- Complete replacement of all residues in the RGD sequence has been described in EP 372 486, assigned to Hoffmann La Roche, where platelet aggregation inhibitors that are derivatives of benzoic and phenylacetic acid are presented. A benzoic acid derivative inhibitor having a particular low IC<sub>50</sub> in an ELISA measurement of fibrinogen GPIIb/IIIa binding has the following structure:



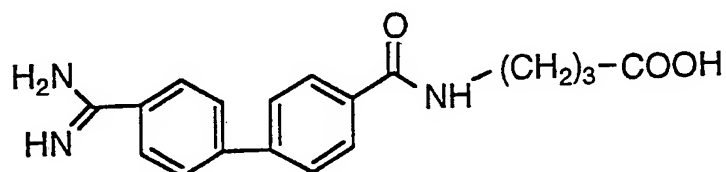
Fibrinogen receptor antagonists possessing similar structures can be found in EP 478 362, EP 478 363, and EP 478 328, all assigned to Merck. A representative Merck compound has the following structure:

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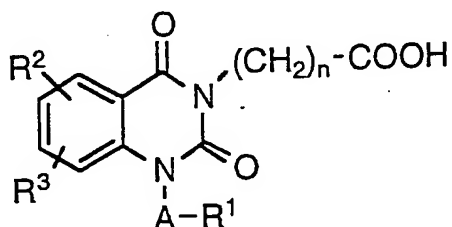
Also of interest are biphenyl derivatives described in EP 483 667 and EP 496 378, the latter publication providing a representative compound having the structure;

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Quinazoline-3-alkanoic acid derivatives are also reported to have an inhibitory effect on platelet aggregation(although possibly by a different mechanism) in EP 456 835 A1. A generic formula representing these compounds is given by:

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where n is 1 to 3 and R<sup>2</sup> and R<sup>3</sup> may be *interalia* hydrogen, lower alkyl, lower alkoxy, aralkyl groups that may be substituted with *interalia* -NR<sup>4</sup>R<sup>5</sup>, where R<sup>4</sup> or R<sup>5</sup> may be hydrogen, lower alkyl, or connected with each other to make five- or six-membered heterocycles which may contain another heteroatom, and A-R<sup>1</sup> may be lower alkyl.

#### 20 D. Benzodiazepines

It is well established that benzodiazepines and related ligands interact with a specific site commonly referred to as the "benzodiazepine receptor" that is associated with a neuro-

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inhibitory postsynaptic GABA receptor and a chloride ionophore channel (see eg. Watjen *et al.*, *J. Med. Chem.* 32:2282-2291[1989]). Binding of ligands to this receptor is known to produce a wide variety of neuro-physiological effects. Benzodiazepines have not been reported to have platelet aggregation inhibition activity. The preparation and therapeutic use of benzodiazepines is described in, for example; EP 0 059 390, EP 0 059 386, and EP 0 394 101.

Benzodiazepinediones have been employed as intermediates in the synthesis of various anti-HIV-1 compounds. For example Kukla, M. J. *et al.*, *J. Med. Chem.* 34:3187-3197 (1991) reduce the dilactam to either the corresponding "one" or diamine in the preparation of various TIBO derivatives having anti-HIV activity. Pyridodiazepines are also described as useful for treatment of HIV infection in U.S. Patent No. 5,087,625.

#### E. Objects

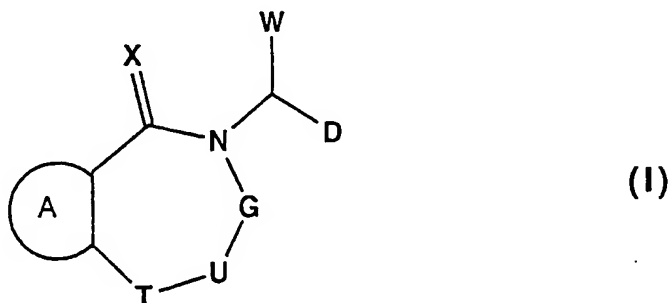
It is an object of this invention to produce nonpeptidyl compounds having potent antithrombotic activity. It is another object of the invention to produce such compounds that are essentially free of amide bonds, substantially rigid, and stable to degradation. It is a further object to produce potent nonpeptidyl antithrombotics that specifically inhibit the GPIIbIIIa-Fg interaction but do not strongly inhibit other RGD sensitive integrin interactions including the Vn-VnR, Fn-FnR, and GPIIbIIIa-vWF interactions. It is still a further object to produce potent nonpeptidyl platelet aggregation inhibitors that do not significantly increase cutaneous bleeding time or diminish other hemodynamic factors. It is also an object of this invention to produce nonpeptidyl compounds having a long half-life and a large therapeutic range.

It is still a further object to produce nonpeptidyl compounds that are capable of inhibiting other integrin interactions such as the Vn-VnR interaction.

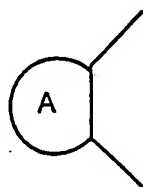
These and other objects of this invention will be apparent from consideration of the invention as a whole.

#### SUMMARY OF THE INVENTION

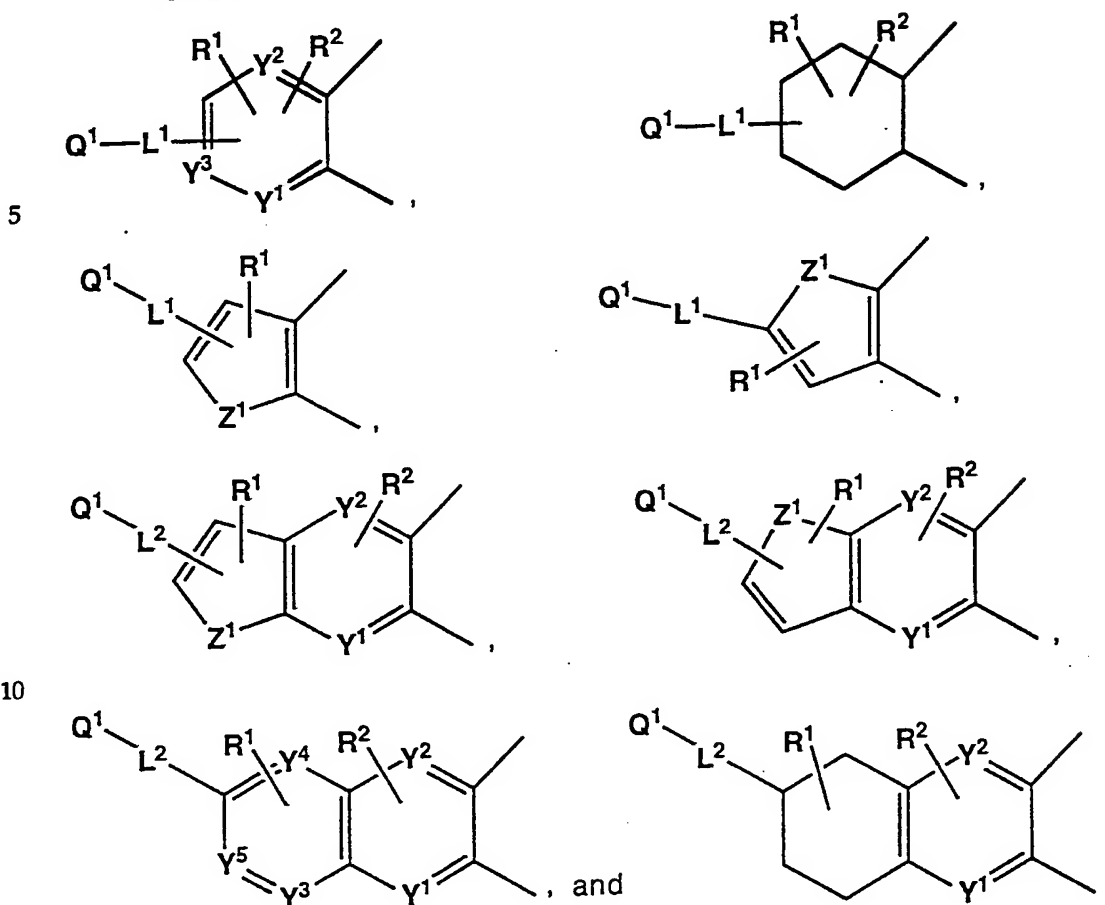
The objects of this invention are accomplished by providing a nonpeptidyl compound represented by structural formula I:



where the partial structure



represents;



where  $R^1$  and  $R^2$  are one to three optional groups typically selected from; hydrogen, halo(F, Cl, Br, I), cyano, carboxamido, carbamoyloxy, formyloxy, formyl, azido, nitro, ureido, thioureido, thiocyanato, hydroxy, mercapto, sulfonamido, and an optionally substituted radical selected from; C<sub>1</sub>-C<sub>12</sub> alkyl, C<sub>2</sub>-C<sub>12</sub> alkenyl, C<sub>3</sub>-C<sub>12</sub> alkynyl, C<sub>3</sub>-C<sub>12</sub> cycloalkyl, C<sub>6</sub>-C<sub>14</sub> aryl, C<sub>6</sub>-C<sub>10</sub> aryl-C<sub>1</sub>-C<sub>8</sub> alkyl, C<sub>1</sub>-C<sub>12</sub> alkyloxy, C<sub>6</sub>-C<sub>14</sub> aryloxy, C<sub>1</sub>-C<sub>12</sub> acylamino, N,N-di(C<sub>1</sub>-C<sub>12</sub>)acylamino, N-(C<sub>1</sub>-C<sub>12</sub>)alkyl-N-(C<sub>1</sub>-C<sub>12</sub>)-acylamino, C<sub>1</sub>-C<sub>12</sub> alkylsulfonamido, N-(C<sub>1</sub>-C<sub>12</sub>)alkyl-N-(C<sub>1</sub>-C<sub>12</sub>)alkyl-sulfonylamino, C<sub>1</sub>-C<sub>12</sub> alkylthiocarbonyl, C<sub>1</sub>-C<sub>12</sub> alkylthio, C<sub>1</sub>-C<sub>12</sub> alkylsulfinyl, C<sub>1</sub>-C<sub>12</sub> alkylsulfonyl, C<sub>1</sub>-C<sub>12</sub> alkylsulfonato, N-(C<sub>1</sub>-C<sub>12</sub>)alkylsulfonamido, N,N-di-(C<sub>1</sub>-C<sub>12</sub>) sulfonamido, N-(C<sub>1</sub>-C<sub>12</sub>) alkyl-N-thioformylamino, C<sub>1</sub>-C<sub>12</sub> thioacylamino, N-(C<sub>1</sub>-C<sub>12</sub>)alkyl-N-(C<sub>1</sub>-C<sub>12</sub>) thioacylamino, C<sub>1</sub>-C<sub>12</sub> alkylsulfinamido, N-(C<sub>1</sub>-C<sub>12</sub>)alkyl-N-(C<sub>1</sub>-C<sub>12</sub>)

C<sub>12</sub>)alkylsulfinylamino, C<sub>1</sub>-C<sub>12</sub> carbalkoxy, C<sub>1</sub>-C<sub>12</sub> alkylcarbonyl, C<sub>1</sub>-C<sub>12</sub> alkanoyloxy, N-(C<sub>1</sub>-C<sub>12</sub>)alkylcarboxamido, N,N-di-(C<sub>1</sub>-C<sub>12</sub>)carboxamido, N-(C<sub>1</sub>-C<sub>12</sub>)alkylcarbamoyloxy, N,N-di-(C<sub>1</sub>-C<sub>12</sub>)carbamoyloxy, and heterocycloalkyl or heteroaryl having from 1 to 3 rings, each ring having from 5 to 7 atoms with from 0-3 heteroatoms selected from N, O, and S, provided that at least one ring contains a heteroatom, where the substituents are typically selected from halo (F, Cl, Br, I), cyano, azido, nitro, hydroxy, mercapto, sulfonamido, ureido, thioureido, carboxamido, carbamoyloxy, formyloxy, formyl, C<sub>1</sub>-C<sub>4</sub>alkyl, C<sub>1</sub>-C<sub>4</sub>alkoxy, phenyl, and phenoxy. Optionally, R<sup>1</sup> and R<sup>2</sup> when bonded to adjacent carbon atoms may join to form an unsubstituted or substituted aryl ring, where the substituents are typically selected from halo (F, Cl, Br, I), cyano, azido, nitro, hydroxy, mercapto, sulfonamido, ureido, thioureido, carboxamido, carbamoyloxy, formyloxy, formyl, C<sub>1</sub>-C<sub>4</sub>alkyl, C<sub>1</sub>-C<sub>4</sub>alkoxy, phenyl, and phenoxy.

Y<sup>1</sup>, Y<sup>2</sup>, Y<sup>3</sup>, Y<sup>4</sup> and Y<sup>5</sup> are independently selected from CH, CR<sup>1</sup>, CR<sup>2</sup>, and N; X is O, S, =NCN, or =NO(C<sub>1</sub>-C<sub>3</sub>)alkyl; and Z<sup>1</sup> is selected from CH<sub>2</sub>, NH, S, and O;

Q<sup>1</sup> is a substituted or unsubstituted positively charged nitrogen-containing moiety. Preferably Q<sup>1</sup> is selected from:

(A) an amino group including;

(1) -NH<sub>2</sub>,

(2) -NR<sup>3</sup>H,

(3) -NR<sup>3</sup>R<sup>4</sup>, and

(4) -NR<sup>3</sup>R<sup>4</sup>R<sup>5</sup>,

where R<sup>3</sup>, R<sup>4</sup>, and R<sup>5</sup> are typically selected from; (i) an optionally substituted radical selected from (a) -NR<sup>6</sup>R<sup>7</sup>, (b) -C(=NR<sup>8</sup>)-NR<sup>6</sup>R<sup>7</sup>, (c) -N=CR<sup>9</sup>-NR<sup>6</sup>R<sup>7</sup>, (d) -NR<sup>10</sup>-CR<sup>9</sup>=NR<sup>8</sup>, and (e) -NR<sup>10</sup>-C(=NR<sup>8</sup>)-NR<sup>6</sup>R<sup>7</sup> where each R<sup>6</sup>, R<sup>7</sup>, R<sup>8</sup>, R<sup>9</sup>, and R<sup>10</sup> is independently selected from hydrogen, C<sub>1</sub>-C<sub>4</sub>alkoxy, C<sub>1</sub>-C<sub>4</sub>alkyl, and halo(F, Cl, Br, I)-C<sub>1</sub>-C<sub>4</sub>alkyl, (ii) optionally substituted C<sub>1</sub>-C<sub>12</sub>alkyl, (iii) optionally substituted C<sub>3</sub>-C<sub>7</sub>alkenyl, (iv) optionally substituted C<sub>3</sub>-C<sub>7</sub>alkynyl, (v) optionally substituted C<sub>3</sub>-C<sub>12</sub>cycloalkyl, (vi) optionally substituted C<sub>5</sub>-C<sub>12</sub>cycloalkenyl, (vii) optionally substituted C<sub>6</sub>-C<sub>14</sub>aryl, (viii) optionally substituted C<sub>1</sub>-C<sub>6</sub>alkyl-C<sub>6</sub>-C<sub>14</sub>-aryl, (ix) optionally substituted C<sub>3</sub>-C<sub>6</sub>alkenyl-C<sub>6</sub>-C<sub>10</sub>aryl, (x) optionally substituted heterocyclyl, (xi) optionally substituted C<sub>1</sub>-C<sub>6</sub>alkyl-heterocyclyl, (xii) optionally substituted C<sub>1</sub>-C<sub>8</sub>alkoxy, (xiii) optionally substituted C<sub>1</sub>-C<sub>8</sub>thioalkoxy, (xiv) optionally substituted C<sub>3</sub>-C<sub>10</sub>alkenoxy, and (xv) optionally substituted C<sub>6</sub>-C<sub>14</sub>aryloxy, where the substituents are usually one to three R<sup>11</sup>, each R<sup>11</sup> typically selected from (a) optionally substituted C<sub>6</sub>-C<sub>12</sub>aryloxy, (b) optionally substituted C<sub>6</sub>-C<sub>12</sub>arylmino, (c) optionally substituted C<sub>6</sub>-C<sub>12</sub>aroxy, (d) optionally substituted C<sub>6</sub>-C<sub>12</sub>arylthio, where the substituents are usually one to three R<sup>12</sup>, each R<sup>12</sup> typically selected from nitro, amino, C<sub>1</sub>-C<sub>8</sub>alkylamino, di-(C<sub>1</sub>-C<sub>8</sub>)alkylamino, amidino, aminomethyleneimino, imino, imino-C<sub>1</sub>-C<sub>4</sub>alkyl, iminomethyleneamino, guanidino, C<sub>6</sub>-C<sub>10</sub>arylmino, C<sub>1</sub>-C<sub>8</sub>acylamino, C<sub>1</sub>-C<sub>4</sub>alkylsulfonamino, azido, cyano, hydroxy, hydroxy-C<sub>1</sub>-C<sub>8</sub>-alkyl, C<sub>1</sub>-C<sub>8</sub>-alkoxy, phenyloxy, C<sub>1</sub>-C<sub>8</sub>alkanoyloxy, C<sub>1</sub>-C<sub>8</sub>alkanoyl, C<sub>6</sub>-C<sub>12</sub>aroxy,

benzamido, phenyl, halo(F, Cl, Br, I), halo-C<sub>1</sub>-C<sub>8</sub>-alkyl, and C<sub>1</sub>-C<sub>8</sub>-alkyl, (e) C<sub>1</sub>-C<sub>8</sub> alkoxy (f) C<sub>1</sub>-C<sub>8</sub>alkthio (g) halo(F, Cl, Br, I), (h) hydroxy, (i) mercapto, (j) C<sub>1</sub>-C<sub>8</sub> alkylcarbonyl, (k) carbamoyl, (l) formyl, (m) formyloxy, (n) carboxy, (o) carb-C<sub>1</sub>-C<sub>8</sub> alkyloxy, (p) C<sub>1</sub>-C<sub>8</sub> alkanoyloxy, (q) N-(C<sub>1</sub>-C<sub>8</sub>)-alkylcarboxamido, (r) N-(C<sub>1</sub>-C<sub>8</sub>), N-(C<sub>1</sub>-C<sub>8</sub>)-dialkylcarboxamido, (s) carbamoyloxy, (t) N-(C<sub>1</sub>-C<sub>8</sub>) alkylcarbamoyloxy, (u) N-(C<sub>1</sub>-C<sub>8</sub>), N-(C<sub>1</sub>-C<sub>8</sub>)-dialkylcarbamoyloxy, (v) C<sub>1</sub>-C<sub>8</sub> alkylsulfinyl, (w) C<sub>1</sub>-C<sub>8</sub> alkylsulfonyl, (x) C<sub>1</sub>-C<sub>8</sub> alkylsulfonato, (y) sulfo, (z) sulfonamido, (aa) N-(C<sub>1</sub>-C<sub>8</sub>) alkylsulfonamido, (ab) N-(C<sub>1</sub>-C<sub>8</sub>), N-(C<sub>1</sub>-C<sub>8</sub>) dialkylsulfonamido, (ac) amino, (ad) C<sub>1</sub>-C<sub>8</sub> alkylamino, (ae) C<sub>1</sub>-C<sub>8</sub> dialkylamino, (af) C<sub>1</sub>-C<sub>8</sub> acylamino, (ag) N-(C<sub>1</sub>-C<sub>8</sub>), N-(C<sub>1</sub>-C<sub>8</sub>)-diacylamino, (ah) N-(C<sub>1</sub>-C<sub>8</sub>)-alkyl-N-(C<sub>1</sub>-C<sub>8</sub>)-acylamino, (ai) formylamino, (aj) ureido, (ak) isothioureido, (al) amino-C<sub>2</sub>-C<sub>8</sub> alkylthio, (am) amino-C<sub>2</sub>-C<sub>8</sub> alkloxy, (an) amidino, (ao) guanidino, (ap) aminomethyleneimino, (aq) imino, (ar) imino-C<sub>1</sub>-C<sub>4</sub> alkyl, (as) iminomethyleneamino, (at) glycyllamino, (au) glycyll, (av) phthalimido, (aw) succinimido, (ax) morpholino, (ay) C<sub>1</sub>-C<sub>8</sub> alkylsulfonamido, (az) N-(C<sub>1</sub>-C<sub>8</sub>)-alkyl-N-(C<sub>1</sub>-C<sub>8</sub>) alkyl sulfonoylamino, (ba) C<sub>1</sub>-C<sub>8</sub> alkylsulfinamino, (bb) N-(C<sub>1</sub>-C<sub>8</sub>) alkyl-N-(C<sub>1</sub>-C<sub>8</sub>) alkylsulfinamino, (bc) C<sub>1</sub>-C<sub>8</sub> alkoxyamino, (bd) C<sub>1</sub>-C<sub>8</sub> alkoxyamino, (be) N-(C<sub>1</sub>-C<sub>8</sub>) alkyl-N-(C<sub>1</sub>-C<sub>8</sub>) alkoxyamino, (bf) C<sub>3</sub>-C<sub>7</sub> cycloalkyl, (bg) oxo, and (bh) heterocyclyl, optionally any one or two pairs of R<sup>3</sup>-R<sup>10</sup> may independently be joined to form one or two optionally substituted heterocyclic rings, each ring optionally fused with one or two optionally substituted homocyclic or heterocyclic rings of from four to seven atoms where any heterocyclic ring contains from one to four heteroatoms selected from N, O, and S and where any ring may be substituted with from one to three R<sup>12</sup>,

(B) an amidino (aminoiminomethyl) group including;

- (1) -C(=NH)-NH<sub>2</sub>,
- (2) -C(=NH)-NHR<sup>3</sup>,
- (3) -C(=NR<sup>4</sup>)-NHR<sup>3</sup>,
- (4) -C(=NH)-NR<sup>3</sup>R<sup>4</sup>, and
- (5) -C(=NR<sup>5</sup>)-NR<sup>3</sup>R<sup>4</sup>,

where R<sup>3</sup>, R<sup>4</sup>, and R<sup>5</sup> are defined above,

(C) an aminoalkyleneamino group including;

- (1) -N=CH-NH<sub>2</sub>,
- (2) -N=CH-NHR<sup>3</sup>,
- (3) -N=CH-NR<sup>3</sup>R<sup>4</sup>, and
- (4) -N=CR<sup>5</sup>-NR<sup>3</sup>R<sup>4</sup>,

where R<sup>3</sup>, R<sup>4</sup>, and R<sup>5</sup> are defined above,

(D) an iminoalkyleneamino group, including;

- (1) -NH-CH=NH,
- (2) -NH-CH=NR<sup>3</sup>,
- (3) -NH-CR<sup>4</sup>=NR<sup>3</sup>, and
- (4) -NR<sup>5</sup>-CR<sup>4</sup>=NR<sup>3</sup>,



where  $R^3$ ,  $R^4$ , and  $R^5$  are defined above,

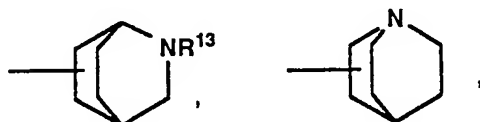
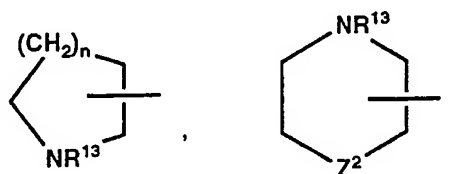
(E) a guanidino (aminoiminomethyleneamino) group including:

- (1)  $-\text{NH}-\text{C}(=\text{NH})-\text{NH}_2$ ,
- (2)  $-\text{NH}-\text{C}(=\text{NH})-\text{NR}^3\text{H}$ ,
- (3)  $-\text{NH}-\text{C}(=\text{NH})-\text{NR}^3\text{R}^4$ ,
- (4)  $-\text{NH}-\text{C}(=\text{NR}^5)-\text{NR}^3\text{R}^4$ ,
- (5)  $-\text{NR}^3-\text{C}(=\text{NR}^3)-\text{NR}^3\text{R}^4$ ,
- (6)  $-\text{NR}^3-\text{C}(=\text{NH})-\text{NR}^3\text{R}^4$ ,
- (7)  $-\text{NR}^3-\text{C}(=\text{NR}^3)-\text{NH}_2$ ,
- (8)  $-\text{NR}^3-\text{C}(=\text{NH})-\text{NH}_2$ ,
- (9)  $-\text{NR}^3-\text{C}(=\text{NR}^3)-\text{NHR}^4$ , and
- (10)  $-\text{NR}^3-\text{C}(=\text{NH})-\text{NHR}^4$ ,

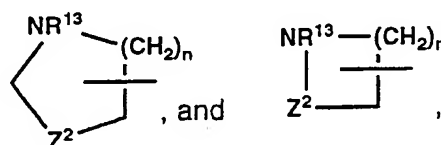
where  $R^3$ ,  $R^4$ , and  $R^5$  are defined above,

(F) an optionally substituted saturated heterocyclic group including:

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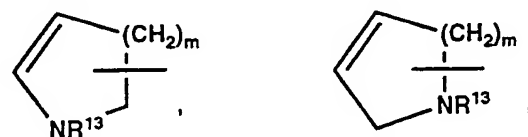
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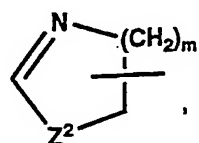
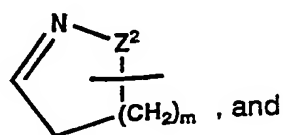
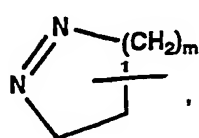


where (1)  $n$  is 0, 1, 2, or 3, (2)  $R^{13}$  is selected from:  $R^6$ ,  $-\text{CR}^9=\text{NR}^8$ ,  $-\text{CR}^9(=\text{NR}^8)-\text{NR}^6\text{R}^7$ ,  $-\text{C}(=\text{NR}^8)-\text{NR}^6\text{R}^7$ ,  $-\text{N}=\text{CR}^9-\text{NR}^6\text{R}^7$ ,  $-\text{NR}^{10}-\text{CR}^9=\text{NR}^8$ , and  $-\text{NR}^{10}-\text{C}(=\text{NR}^8)-\text{NR}^6\text{R}^7$  where  $R^6$ - $R^{10}$  are defined above, (3)  $Z^2$  is O, S, or  $\text{NR}^{13}$ , and (4) the substituents are independently one to three  $R^{12}$ ,

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(G) an optionally substituted unsaturated (nonaromatic) heterocyclyl including:

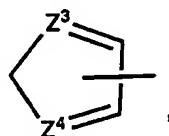
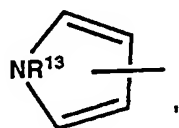




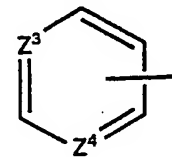
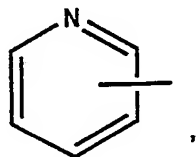
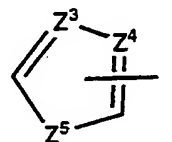
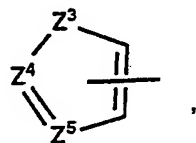
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where (1)  $m$  is 1, 2, or 3, (2)  $Z^2$  and  $R^{13}$  are defined above, and (3) the substituents are typically one to three  $R^{12}$ ,

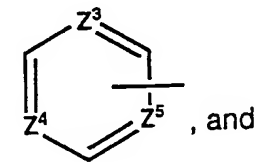
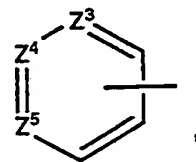
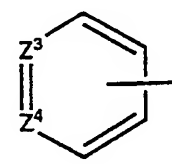
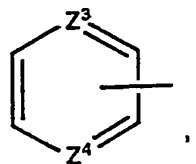
(H) an optionally substituted unsaturated (aromatic) heterocyclyl including:



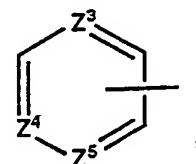
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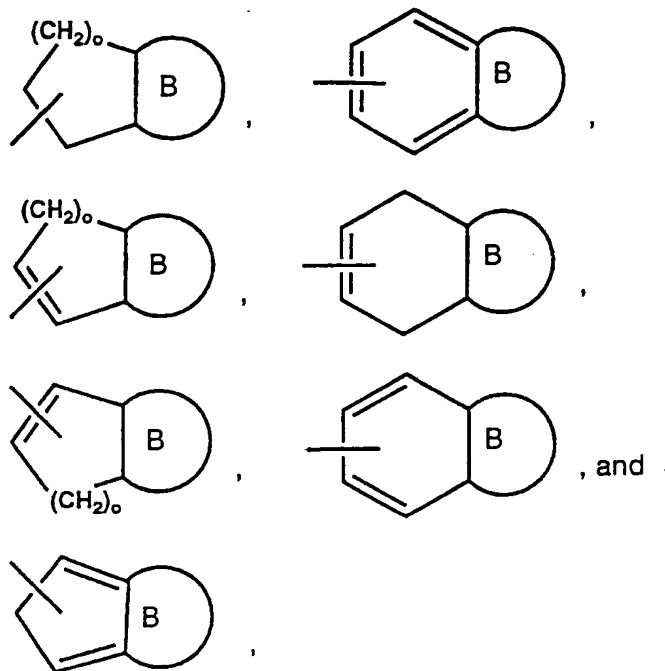
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where (1)  $Z^3$ ,  $Z^4$ , and  $Z^5$  are typically selected from O, S, N, and NH, provided at least one  $Z^3$ ,  $Z^4$ , or  $Z^5$  is N or NH, (2)  $R^{13}$  is defined above, and the substituents are independently one to three  $R^{12}$ ,

(I) an optionally substituted bicycloheterocyclic group including:

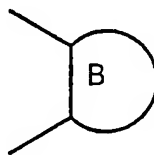
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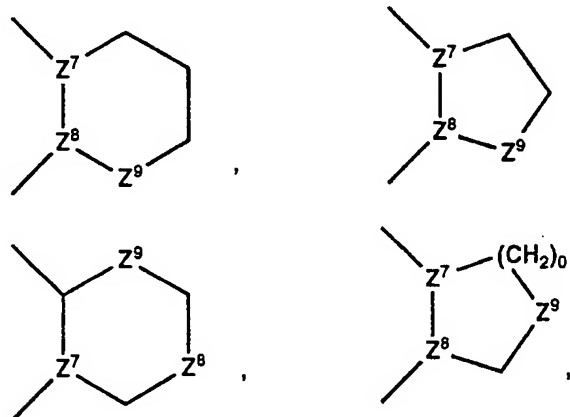
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where the partial structure



represents



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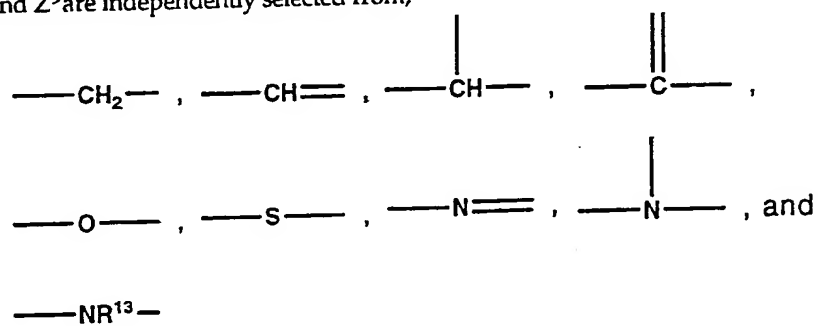
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where  $Z^7$ ,  $Z^8$ , and  $Z^9$  are independently selected from;

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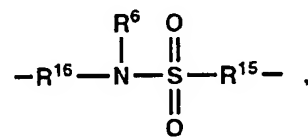
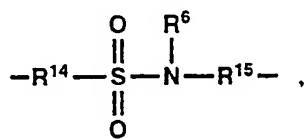
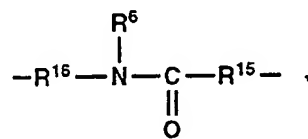
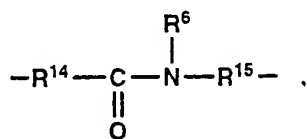
provided that at least one  $Z^7$ ,  $Z^8$ , or  $Z^9$  is



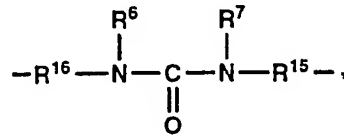
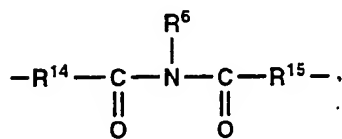
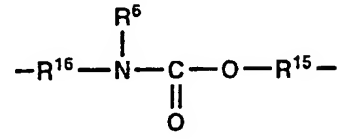
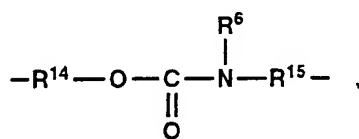
where (1) o is 0, 1, or 2, (2) R<sup>13</sup> is defined above, and (3) the substituents are independently one to three R<sup>12</sup>. Exemplary heterocyclic Q<sup>1</sup> groups include; isoindoliny, quinuclidiny, morpholiny, and 1,3-diazacyclohex-4-ene. Optionally, any of the nitrogen containing heterocycles described above may be substituted with amino, imino, amidino, aminomethyleneimino, iminomethyleneamino, guanidino, N<sup>G</sup>-aminoguanidino, alkylamino, dialkylamino, trialkylamino, or alkylideneamino groups.

L<sup>1</sup> is a bivalent radical containing from 3 to 9 methylene groups where any methylene group or groups may be replaced with one or more alkene, alkyne, aryl, or functional groups containing the heteroatoms selected from the group N, O, and S. Preferably, L<sup>1</sup> is a saturated or unsaturated, linear, branched, or cyclic bivalent radical separating Q from the 7 carbon of the benzodiazepinedione nucleus. Typically, the heteroatoms (N, O, and S) will comprise from 0-5 of the atoms separating Q from the benzodiazepinedione nucleus. Preferably, L<sup>1</sup> is typically an optionally substituted bivalent radical including: (A) C<sub>3</sub>-C<sub>7</sub>-alkylene, (B) C<sub>3</sub>-C<sub>7</sub>-cycloalkylene, (C) C<sub>3</sub>-C<sub>7</sub>-alkenylene, (D) C<sub>3</sub>-C<sub>7</sub>-alkadienylene, (E) C<sub>3</sub>-C<sub>7</sub>-alkynylene, (F) C<sub>4</sub>-C<sub>7</sub>-alkadiynylene, (G) C<sub>4</sub>-C<sub>7</sub>-alkenynylene, (H) C<sub>6</sub>-C<sub>14</sub>-arylene, (I) C<sub>6</sub>-C<sub>14</sub>-aryl-C<sub>2</sub>-C<sub>4</sub>-alkynylene, (J) C<sub>1</sub>-C<sub>3</sub>-alkyl-C<sub>6</sub>-C<sub>14</sub>-aryl-C<sub>2</sub>-C<sub>4</sub>-alkynylene, (K) C<sub>6</sub>-C<sub>14</sub>-aryl-C<sub>2</sub>-C<sub>4</sub>-alkenylene, (L) C<sub>1</sub>-C<sub>3</sub>-alkyl-C<sub>6</sub>-C<sub>14</sub>-arylene, (M) C<sub>1</sub>-C<sub>3</sub>-alkyl-C<sub>6</sub>-C<sub>14</sub>-aryl-C<sub>2</sub>-C<sub>4</sub>-alkenylene, (N) C<sub>6</sub>-C<sub>14</sub>-aryl-C<sub>1</sub>-C<sub>3</sub>-alkylene, (O) C<sub>6</sub>-C<sub>14</sub>-aryl-C<sub>1</sub>-C<sub>3</sub>-alkyloxyene, (P) C<sub>1</sub>-C<sub>2</sub>-alkyl-C<sub>6</sub>-C<sub>14</sub>-aryl-C<sub>1</sub>-C<sub>2</sub>-alkylene, (Q) C<sub>1</sub>-C<sub>3</sub>-alkyloxy-C<sub>6</sub>-C<sub>14</sub>-arylene, (R) C<sub>2</sub>-C<sub>6</sub>-alkyloxyene, (S) C<sub>1</sub>-C<sub>5</sub>-alkyloxy-C<sub>1</sub>-C<sub>5</sub>-alkylene, (T) C<sub>6</sub>-C<sub>10</sub>-aryloxyene, (U) C<sub>6</sub>-C<sub>10</sub>-aryloxy-C<sub>1</sub>-C<sub>5</sub>-alkylene, (V) C<sub>2</sub>-C<sub>6</sub>-alkylthioene, (W) C<sub>1</sub>-C<sub>5</sub>-alkylthio-C<sub>1</sub>-C<sub>5</sub>-alkylene, (X) C<sub>6</sub>-C<sub>10</sub>-arylthioene, (Y) C<sub>6</sub>-C<sub>10</sub>-arylthio-C<sub>1</sub>-C<sub>5</sub>-alkylene, (Z) C<sub>1</sub>-C<sub>5</sub>-alkylsulfoxide-C<sub>1</sub>-C<sub>5</sub>-alkylene, (AA) C<sub>1</sub>-C<sub>5</sub>-alkylsulfone-C<sub>1</sub>-C<sub>5</sub>-alkylene,

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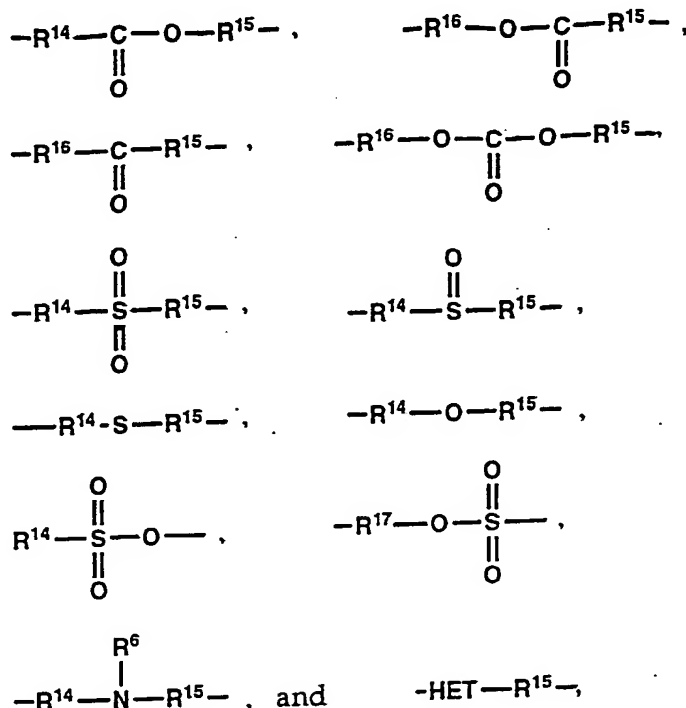
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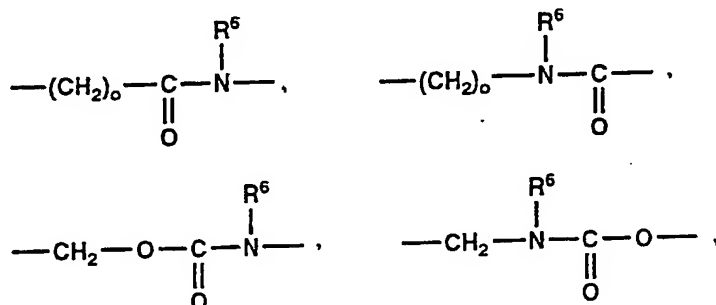
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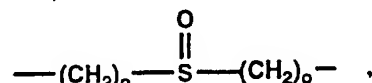
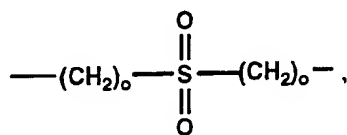
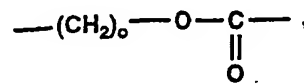
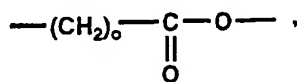
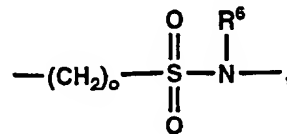
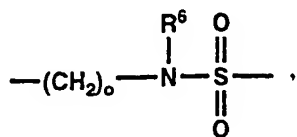
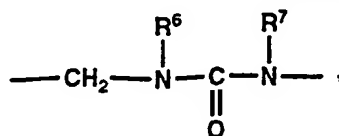
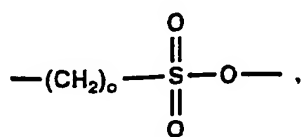


where (1)  $R^{14}$  is selected from; (i) a chemical bond, (ii)  $C_1$ - $C_5$ -alkyl, (iii)  $C_3$ - $C_7$ -cycloalkyl, (iv)  $C_2$ - $C_5$ -alkenyl, (v)  $C_3$ - $C_5$ -alkynyl, (vi)  $C_6$ - $C_{10}$ -aryl, (vii)  $C_1$ - $C_2$ -alkyl- $C_6$ - $C_{10}$ -aryl- $C_1$ - $C_2$ -alkyl, (viii)  $C_1$ - $C_3$ -alkyl- $C_6$ - $C_{12}$ -aryl, (ix)  $C_6$ - $C_{10}$ -aryl- $C_1$ - $C_2$ -alkyl, (x)  $C_6$ - $C_{10}$ -aryloxy- $C_1$ - $C_2$ -alkyl, and (xi) piperiziny, (2)  $R^{15}$  is selected from; (i) a chemical bond, (ii)  $C_1$ - $C_4$ -alkyl, (iii)  $C_2$ - $C_4$ -alkenyl, (iv)  $C_2$ - $C_4$ -alkynyl, (v)  $C_1$ - $C_3$ -alkyl- $C_6$ - $C_{12}$ -aryl, and (vi)  $C_6$ - $C_{10}$ -aryl, (3)  $R^{16}$  is selected from; (i) a chemical bond, (ii)  $C_1$ - $C_5$ -alkyl, (iii)  $C_3$ - $C_7$ -cycloalkyl, (iv)  $C_3$ - $C_5$ -alkenyl, (v)  $C_3$ - $C_5$ -alkynyl, (vi)  $C_6$ - $C_{10}$ -aryl, (vii)  $C_1$ - $C_3$ -alkyl- $C_6$ - $C_{12}$ -aryl, (viii)  $C_6$ - $C_{10}$ -aryl- $C_1$ - $C_2$ -alkyl, and (ix) piperiziny, (4)  $R^{17}$  is selected from; (i)  $C_3$ - $C_4$ -alkenyl, (ii)  $C_3$ - $C_4$ -alkynyl, (iii)  $C_6$ - $C_{10}$ -aryl, and (iv) benzyl, (5) HET is a saturated or unsaturated heterocycle having from 5-14 atoms in the cycle(s) and from 1-3 heteroatoms selected from N, O, and S, where the substituents are selected from one to three  $R^{12}$ .

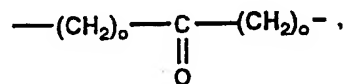
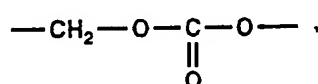
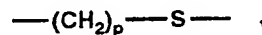
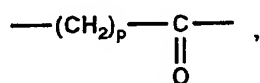
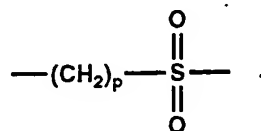
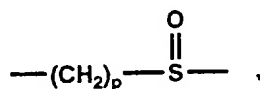
$L^2$  is typically an optionally substituted bivalent radical including but not limited to;



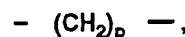
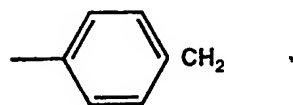
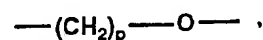
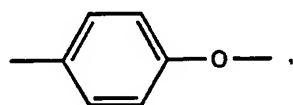
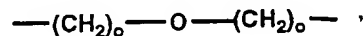
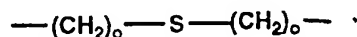
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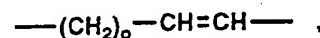
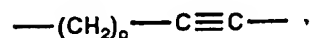
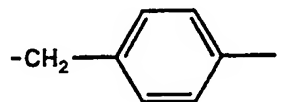
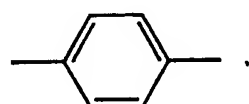
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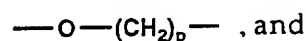
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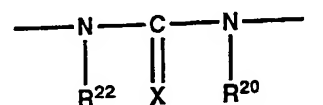
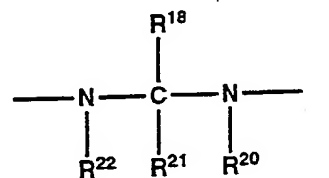
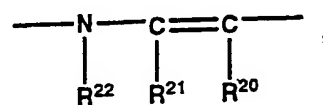
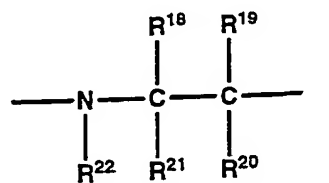
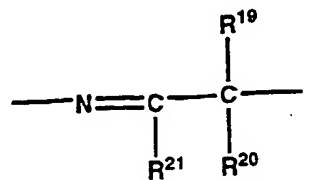
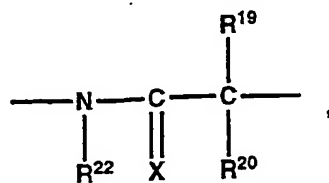
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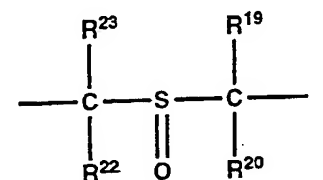
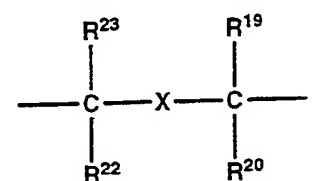
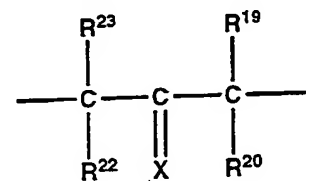
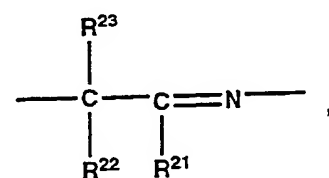
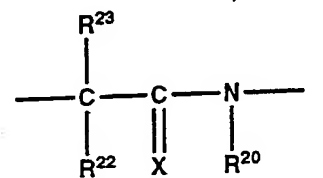
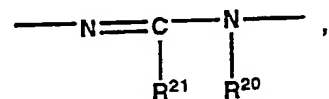
where; (1) p is 1, 2, 3 or 4, (2) R<sup>6</sup>, R<sup>7</sup>, o, and HET are defined above, and (3) the substituents are typically selected from one to three R<sup>12</sup>.

T-U-G is selected from;

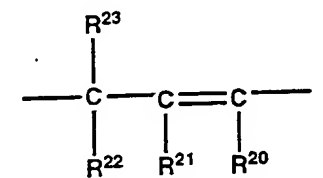
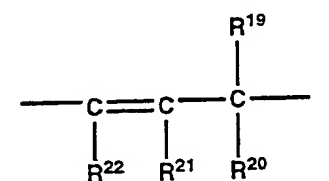
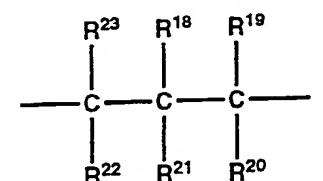
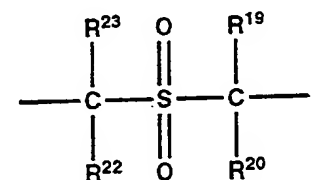
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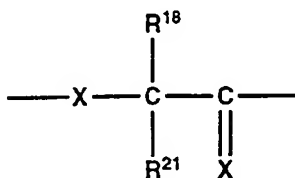
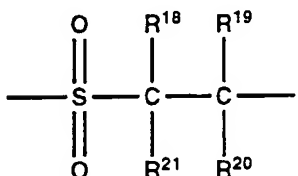
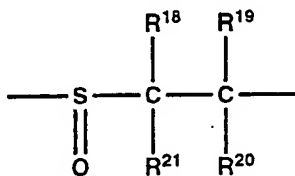
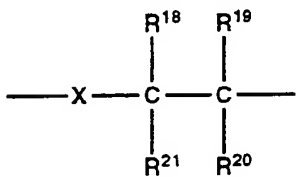
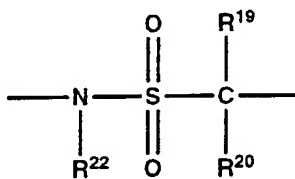
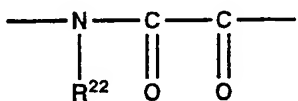
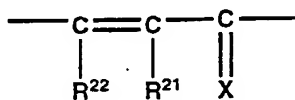
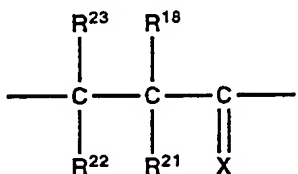
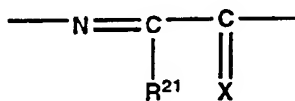
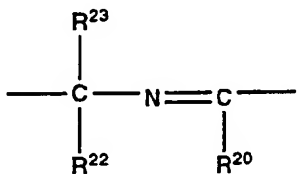
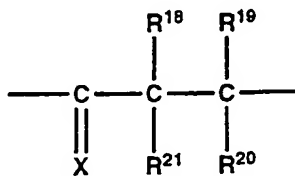
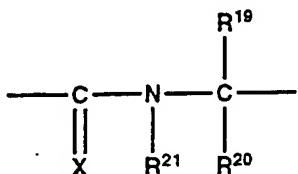
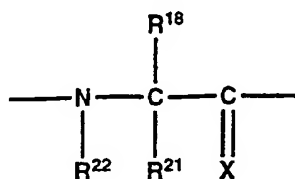
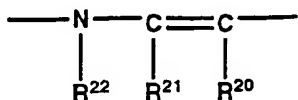
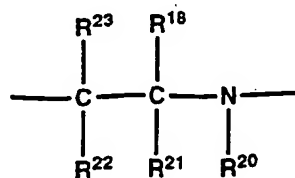
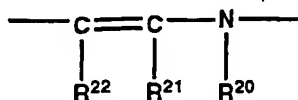
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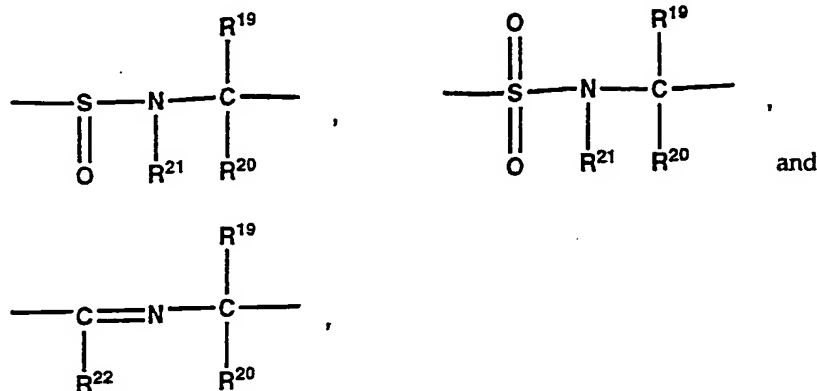


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- 5 where; (1) X is defined above, (2) R<sup>18</sup>, R<sup>19</sup>, R<sup>20</sup>, R<sup>23</sup>, are typically selected from (i) hydrogen, (ii) optionally substituted C<sub>1</sub>-C<sub>12</sub> alkyl, (iii) optionally substituted C<sub>3</sub>-C<sub>12</sub> alkenyl, (iv) optionally substituted C<sub>3</sub>-C<sub>14</sub> cycloalkyl, (v) optionally substituted C<sub>1</sub>-C<sub>12</sub> alkyl-C<sub>6</sub>-C<sub>14</sub> aryl, (vi) optionally substituted C<sub>6</sub>-C<sub>14</sub> aryl, (vii) optionally substituted C<sub>1</sub>-C<sub>4</sub> alkylphenyl, and (viii) optionally substituted C<sub>1</sub>-C<sub>12</sub> alkoxy, (3) R<sup>21</sup> and R<sup>22</sup> are typically selected from (i) Q<sup>2</sup>-
- 10 L<sup>3</sup>- where Q<sup>2</sup> is selected from hydrogen and Q<sup>1</sup>, and L<sup>3</sup> is either a chemical bond, L<sup>1</sup> or L<sup>2</sup>, (ii) optionally substituted C<sub>1</sub>-C<sub>12</sub>-alkyl, (iii) optionally substituted C<sub>6</sub>-C<sub>14</sub>-aryl, (iv) optionally substituted C<sub>3</sub>-C<sub>14</sub>-cycloalkyl, (v) optionally substituted C<sub>1</sub>-C<sub>12</sub>-alkyl-C<sub>6</sub>-C<sub>14</sub> aryl, and (vi) optionally substituted C<sub>1</sub>-C<sub>12</sub>-alkyl-C<sub>3</sub>-C<sub>14</sub>-cycloalkyl, where the substituents are usually (a) halo (F, Cl, Br, I), (b) nitro, (c) hydroxy, (d) carboxy, (e) tetrazole, (f) hydroxamate, (g)
- 15 sulfonamide, (h) trifluoroimide, (i) phosphonate, (j) C<sub>1</sub>-C<sub>6</sub>-alkyl, (k) C<sub>6</sub>-C<sub>14</sub>-aryl, (l) benzyl, (m) C<sub>3</sub>-C<sub>14</sub>-cycloalkyl, (n) COR<sup>24</sup> where R<sup>24</sup> is selected from C<sub>1</sub>-C<sub>8</sub>-alkoxy, C<sub>3</sub>-C<sub>12</sub>-alkenoxy, C<sub>6</sub>-C<sub>12</sub>-aryloxy, di-C<sub>1</sub>-C<sub>8</sub>-alkylamino-C<sub>1</sub>-C<sub>8</sub>-alkoxy, acylamino-C<sub>1</sub>-C<sub>8</sub>-alkoxy selected from acetylaminooethoxy, nicotinoylaminoethoxy, succinamidooethoxy, and pivaloyloxyethoxy, and C<sub>6</sub>-C<sub>12</sub>-aryl-C<sub>1</sub>-C<sub>8</sub>-alkoxy where the aryl group is unsubstituted or substituted with one to
- 20 three of the groups nitro, halo (F, Cl, Br, I), C<sub>1</sub>-C<sub>4</sub>-alkoxy, amino, hydroxy, hydroxy-C<sub>2</sub>-C<sub>8</sub>-alkoxy, and dihydroxy-C<sub>3</sub>-C<sub>8</sub>-alkoxy, and (o) CONR<sup>25</sup>R<sup>26</sup> where R<sup>25</sup> and R<sup>26</sup> are typically selected from hydrogen, C<sub>1</sub>-C<sub>10</sub>-alkyl, C<sub>3</sub>-C<sub>10</sub>-alkenyl, C<sub>6</sub>-C<sub>14</sub>-aryl, C<sub>1</sub>-C<sub>6</sub>-alkyl-C<sub>6</sub>-C<sub>10</sub>-aryl. Optionally, R<sup>25</sup> and R<sup>26</sup> taken together may form trimethylene, tetramethylene, pentamethylene, and 3-oxopentamethylene. Also, optionally, R<sup>1</sup> and R<sup>22</sup> taken together may
- 25 form an optionally substituted 5, 6, or 7 member saturated or unsaturated homocyclic or heterocyclic ring containing 0-3 heteroatoms selected from O, S, and N, where the substituents are selected from R<sup>21</sup>.

30 D is typically (A) R<sup>21</sup>, or (B) -(C=O)-Xaa, where Xaa is one to three D or L α-amino acid residues.

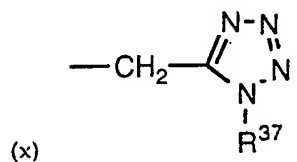
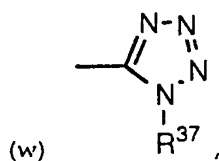
W is -R<sup>27</sup>-w.

R<sup>27</sup> is selected from (a) a covalent bond, (b) substituted or unsubstituted methylene, and (c) substituted or unsubstituted ethylene where the substituents are selected from (i) nitro,

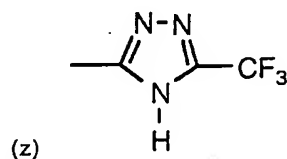
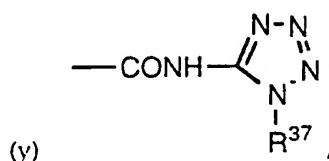
(ii) halo(F, Cl, Br, I), (iii) C<sub>1</sub>-C<sub>6</sub> alkyl, (iv) halo(F, Cl, Br, I)-C<sub>1</sub>-C<sub>6</sub> alkyl, and (v) substituted or unsubstituted phenyl where the substituents are selected from (1) C<sub>1</sub>-C<sub>6</sub> alkyl, (2) C<sub>1</sub>-C<sub>6</sub> alkoxy, (3) halo(F, Cl, Br, I), and (4) CF<sub>3</sub>.

- "w" is selected from (a) -COR<sup>28</sup>, (b) -SO<sub>3</sub>R<sup>31</sup>, (c) -NHSO<sub>2</sub>R<sup>32</sup>, (d) -PO(OR<sup>31</sup>)<sub>2</sub>, (e) -SO<sub>2</sub>NHR<sup>32</sup>, (f) -CONHOR<sup>31</sup>, (g) -C(OH)R<sup>33</sup>PO(OR<sup>33</sup>)<sub>2</sub>, (h) -CN, (i) -SO<sub>2</sub>NH-heteroaryl where the heteroaryl is a 5- or 6-member aromatic ring containing 1 to 3 heteroatoms selected from O, N, and S and where the heteroaryl is unsubstituted or substituted with one or two substituents selected from the group (i) -OH, (ii) -SH, (iii) -(C<sub>1</sub>-C<sub>4</sub>alkyl), (iv) -C<sub>1</sub>-C<sub>4</sub>alkoxyl, (v) CF<sub>3</sub>, (vi) halo(F, Cl, Br, I), (vii) NO<sub>2</sub>, (viii) -COOH, (ix) -COO-(C<sub>1</sub>-C<sub>4</sub>alkyl), (x) -NH<sub>2</sub>, (xi) -NH(C<sub>1</sub>-C<sub>4</sub>alkyl), or (xii) -N(C<sub>1</sub>-C<sub>4</sub>alkyl)<sub>2</sub>, (j) -CH<sub>2</sub>SO<sub>2</sub>NH-heteroaryl, (k) -SO<sub>2</sub>NHCOR<sup>33</sup>, (l) -CH<sub>2</sub>SO<sub>2</sub>NHCOR<sup>32</sup>, (m) -CONHSO<sub>2</sub>R<sup>33</sup>, (n) -CH<sub>2</sub>CONHSO<sub>2</sub>R<sup>33</sup>, (o) -NHCONHSO<sub>2</sub>R<sup>33</sup>, (p) -NHSO<sub>2</sub>NHCOR<sup>33</sup>, (q) -CONHNHSO<sub>2</sub>CF<sub>3</sub>, (r) CON(OH)R<sup>31</sup>, (s) -CONHCOCF<sub>3</sub>, (t) -CONHSO<sub>2</sub>R<sup>28</sup>, (u) -CONHSO<sub>2</sub>R<sup>29</sup>, (v) -CONHSO<sub>2</sub>R<sup>30</sup>,

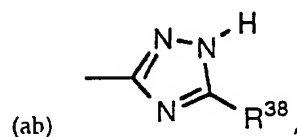
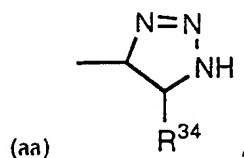
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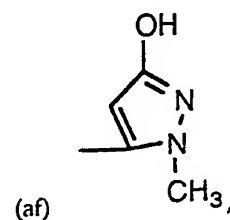
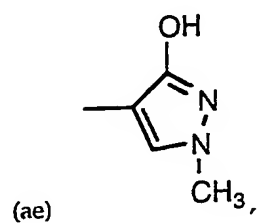
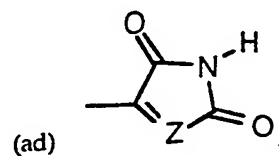
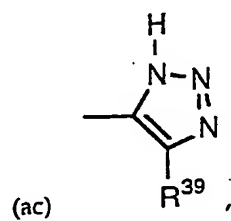


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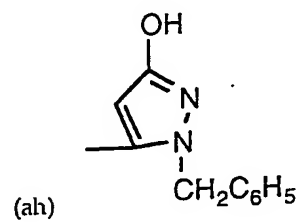
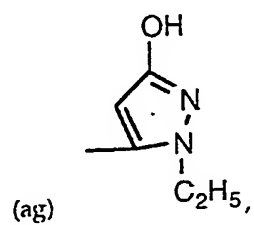


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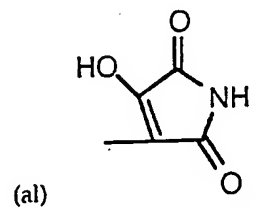
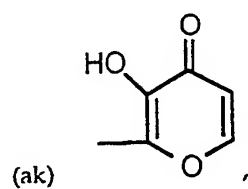
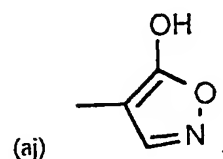
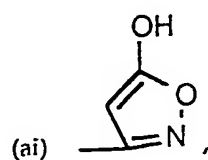




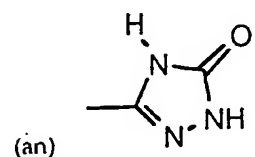
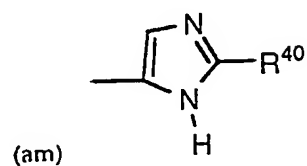
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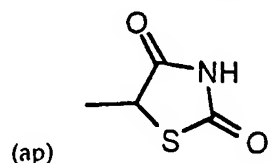
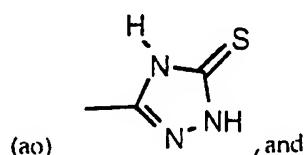


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$R^{28}$  is selected from the group consisting of (a) hydroxy, (b) C<sub>1</sub>-C<sub>8</sub>-alkoxy, (c) C<sub>3</sub>-C<sub>12</sub>-alkenoxy, (d) C<sub>6</sub>-C<sub>12</sub>-aryloxy, (e) C<sub>1</sub>-C<sub>6</sub>-alkyl-C<sub>6</sub>-C<sub>12</sub>-aryloxy, (f) di-C<sub>1</sub>-C<sub>8</sub>-alkylamino-C<sub>1</sub>-C<sub>8</sub>-alkoxy, (g) acylamino-C<sub>1</sub>-C<sub>8</sub>-alkoxy selected from the group (i) acetylaminoethoxy, (ii) nicotinoylaminoethoxy, and (iii) succinamidoethoxy, (h) C<sub>1</sub>-C<sub>8</sub>-alkoxyloxy-C<sub>1</sub>-C<sub>8</sub>-alkoxy, (i) C<sub>6</sub>-C<sub>12</sub>-aryl-C<sub>1</sub>-C<sub>8</sub>-alkoxy where the aryl group is unsubstituted or substituted with one to three of the groups (i) nitro, (ii) halo (F, Cl, Br, I), (iii) C<sub>1</sub>-C<sub>4</sub>-alkoxy, and (iv) amino, (j) hydroxy-C<sub>2</sub>-C<sub>8</sub>-alkoxy, (k) dihydroxy-C<sub>3</sub>-C<sub>8</sub>-alkoxy, and (l)  $NR^{29}R^{30}$ .

$R^{29}$  and  $R^{30}$  are independently selected from the group consisting of (a) hydrogen, (b) C<sub>1</sub>-C<sub>8</sub>-alkyl, (c) C<sub>3</sub>-C<sub>8</sub>-alkenyl, (d) C<sub>6</sub>-C<sub>12</sub>-aryl where the aryl group is unsubstituted or substituted with one to three of the groups (i) nitro, (ii) halo (F, Cl, Br, I), (iii) C<sub>1</sub>-C<sub>4</sub>-alkoxy, and (iv) amino, and (e) C<sub>6</sub>-C<sub>12</sub>-aryl-C<sub>1</sub>-C<sub>8</sub>-alkyl where the aryl group is unsubstituted or substituted with one to three of the groups (i) nitro, (ii) halo (F, Cl, Br, I), and (iii) C<sub>1</sub>-C<sub>4</sub>-alkoxy.

$R^{31}$  is selected from the group consisting of (a) H, (b) C<sub>1</sub>-C<sub>6</sub> alkyl, (c) halo(F, Cl, Br, I)-C<sub>1</sub>-C<sub>6</sub> alkyl, (d) phenyl, (e) benzyl, and (f) -CH<sub>2</sub>-O-COCH<sub>3</sub>.

$R^{32}$  is selected from the group consisting of (a) H, (b) benzyl and (c) -CH( $R^{35}$ )-O-C(O) $R^{35}$ .

$R^{33}$  is selected from the group consisting of (a) aryl, (b) heteroaryl, (c) (C<sub>3</sub>-C<sub>7</sub>-cycloalkyl, (d) (C<sub>1</sub>-C<sub>4</sub>)-alkyl, unsubstituted or substituted with a substituent selected from the group consisting of (i) aryl, (ii) heteroaryl, (iii) -OH, (iv) -SH, (v) (C<sub>1</sub>-C<sub>4</sub>)-alkyl, (vi) (C<sub>1</sub>-C<sub>4</sub>)-alkoxy, (vii) (C<sub>1</sub>-C<sub>4</sub>)-alkylthio, (viii) -CF<sub>3</sub>, (ix) halo (F, Cl, Br, I), (x) -NO<sub>2</sub>, (xi) -CO<sub>2</sub>H, (xii) CO<sub>2</sub>-(C<sub>1</sub>-C<sub>4</sub>)-alkyl, (xiii) -NH<sub>2</sub>, (xiv) -N[(C<sub>1</sub>-C<sub>4</sub>)-alkyl]<sub>2</sub>, (xv) -NH[(C<sub>1</sub>-C<sub>4</sub>)-alkyl], (xvi) -PO<sub>3</sub>H or (xvii) PO(OH)(C<sub>1</sub>-C<sub>4</sub>)-alkoxy, or (e) (C<sub>1</sub>-C<sub>4</sub>)-perfluoroalkyl.

$R^{34}$  is selected from the group consisting of (a) -CN, (b) -NO<sub>2</sub>, (c) -COOR<sup>31</sup>, (d) C<sub>1</sub>-C<sub>6</sub>-perfluoroalkyl, and (e) CF<sub>3</sub>.

$R^{35}$  is independently selected from the group consisting of (a) H, (b) (C<sub>1</sub>-C<sub>6</sub>)-alkyl, (C<sub>2</sub>-C<sub>6</sub>)-alkenyl, (C<sub>2</sub>-C<sub>6</sub>)-alkynyl or (C<sub>3</sub>-C<sub>8</sub>)-cycloalkyl, each of which is unsubstituted or substituted with: (i) OH, (ii) (C<sub>1</sub>-C<sub>4</sub>)-alkoxy, (iii) CO<sub>2</sub>R<sup>33</sup>, (iv) OCOR<sup>33</sup>, (v) CONHR<sup>33</sup>, (vi) CON( $R^{33}$ )<sub>2</sub>, (vii) N( $R^{33}$ )C(O) $R^{33}$ , (viii) NH<sub>2</sub>, (ix) (C<sub>1</sub>-C<sub>4</sub>)-alkylamino, (x) di[(C<sub>1</sub>-C<sub>4</sub>)-alkyl]amino, (xi) aryl, (xii) heteroaryl, (c) -C(O)-aryl, (d) -NO<sub>2</sub>, (e) halo(Cl, Br, I, F), (f) -OH, (g) -OR<sup>36</sup>, (h) (C<sub>1</sub>-C<sub>4</sub>)-perfluoroalkyl, (i) -SH, (j) -S(O)<sub>1-2</sub> (C<sub>1</sub>-C<sub>4</sub>)-alkyl, (k) CO<sub>2</sub>R<sup>33</sup>, (l) -SO<sub>3</sub>H, (m) -NR<sup>33</sup>R<sup>36</sup>, (n) -NR<sup>33</sup>C(O) $R^{36}$ , (o) -NR<sup>33</sup>COOR<sup>32</sup>, (p) -SO<sub>2</sub>NHR<sup>32</sup>, (q) -SO<sub>2</sub>NR<sup>33</sup>R<sup>33</sup>, (r) -NHSO<sub>2</sub>R<sup>32</sup>, (s) -C(O)NHSO<sub>2</sub>R<sup>32</sup>, (t) aryl, (u) heteroaryl, (v) morpholin-4-yl, (w) CONH<sub>2</sub>, or (y) 1H-tetrazol-5-yl.

R<sup>36</sup> is selected from the group consisting of (a) H or (b) (C<sub>1</sub>-C<sub>4</sub>)-alkyl unsubstituted or substituted with (i) NH<sub>2</sub>, (ii) NH[(C<sub>1</sub>-C<sub>4</sub>)-alkyl], (iii) N[(C<sub>1</sub>-C<sub>4</sub>)-alkyl]<sub>2</sub>, (iv) CO<sub>2</sub>H, (v) CO<sub>2</sub>(C<sub>1</sub>-C<sub>4</sub>)-alkyl, (vi) OH, (vii) SO<sub>3</sub>H, or (viii) SO<sub>2</sub>NH<sub>2</sub>.

5 R<sup>37</sup> is selected from the group consisting of (a) H, (b) (C<sub>1</sub>-C<sub>6</sub>)-alkyl, (c) (C<sub>2</sub>-C<sub>6</sub>)-alkenyl, (d) (C<sub>1</sub>-C<sub>6</sub>)-alkoxyalkyl, (e) -CH<sub>2</sub>-O-COCH<sub>3</sub>, or (f) -CH<sub>2</sub>-phenyl, where the phenyl is unsubstituted or substituted with a substituent selected from -NO<sub>2</sub>, -NH<sub>2</sub>, -OH, or -OCH<sub>3</sub>.

R<sup>38</sup>, R<sup>39</sup>, and R<sup>40</sup> are each independently selected from H, Cl, CN, NO<sub>2</sub>, CF<sub>3</sub>, C<sub>2</sub>F<sub>5</sub>, C<sub>3</sub>F<sub>7</sub>, CHF<sub>2</sub>, CH<sub>2</sub>F, CO<sub>2</sub>CH<sub>3</sub>, CO<sub>2</sub>C<sub>2</sub>H<sub>5</sub>, SO<sub>2</sub>CH<sub>3</sub>, SO<sub>2</sub>CF<sub>3</sub> and SO<sub>2</sub>C<sub>6</sub>F<sub>5</sub>, where Z is selected from O, S, NR<sup>41</sup> and CH<sub>2</sub>.

10 R<sup>41</sup> is selected from hydrogen, CH<sub>3</sub>, and CH<sub>2</sub>C<sub>6</sub>H<sub>5</sub> or a tautomer thereof.

The invention further provides a pharmaceutical composition comprising a pharmaceutically acceptable excipient and/or pharmaceutically acceptable hydrates, solvates, and salts of I.

15 The invention provides a method for inhibiting platelet aggregation or reducing platelet aggregation in a mammal comprising administering a platelet aggregation inhibiting amount of the pharmaceutical composition of I. Optionally, the method includes administering the pharmaceutical composition in combination with a peptidyl antithrombotic and/or a thrombolytic agent and/or an anticoagulant.

20

## DETAILED DESCRIPTION OF THE INVENTION

### A. Definitions

Terms used in the claims and specification are defined as set forth below unless otherwise specified.

25 As used herein, the term "positively charged" when used to describe the moiety Q (either Q<sup>1</sup> or Q<sup>2</sup>) means at least 10% of the Q groups are positively charged at physiological pH. This will normally mean that the pK<sub>b</sub> of the nitrogen containing moiety will be about 6.9 or higher. The phrase "positively charged Q moiety;" as used to define Q, is intended to embrace chemical groups which, when attached to the linking moiety L of Formula I, confers basic character to the compound of Formula I. "Basic character" means proton accepting or  
30 electron donating capability, that is, the capacity of the compound of Formula I to be a proton acceptor in the presence of a proton donor, such as water. Optionally, Q will be a prodrug that is capable of being converted, *in vivo*, into a positively charged basic moiety. An example of a basic group having "basic character" is the guanidino (H<sub>2</sub>N-C(=NH)-NH-) group. There are many examples of other equivalent basic groups other than the guanidino group that can be  
35 substituted into Q to produce biologically equivalent compounds of Formula I. Such other basic groups are collectively referred to as "bioisosteres of the guanidino group" or "basic bioisosteres". Representative examples of such "basic bioisosteres" are set forth above under Q (A) through (I).

The term "alkyl" means a branched or unbranched, saturated aliphatic hydrocarbon radical, having the number of carbon atoms specified, or if no number is specified, having up to 12 carbon atoms. Examples of alkyl radicals include methyl, ethyl, n-propyl, isopropyl, n-butyl, iso-butyl, sec-butyl, tert-butyl, n-pentyl, 2-methylbutyl, 2,2-dimethylpropyl, n-hexyl, 2-methylpentyl, 2,2-dimethylbutyl, n-heptyl, 2-methylhexyl, and the like. The terms "lower alkyl", "C<sub>1</sub>-C<sub>6</sub> alkyl" and "alkyl of 1 to 6 carbon atoms" are synonymous and used interchangeably. A preferred "C<sub>1</sub>-C<sub>6</sub> alkyl" group is methyl.

The term "substituted C<sub>n</sub>-C<sub>m</sub> alkyl" where m and n are integers identifying the range of carbon atoms contained in the alkyl group, denotes the above alkyl groups that are substituted by one, two or three halogen, hydroxy, protected hydroxy, amino, protected amino, C<sub>1</sub>-C<sub>7</sub> acyloxy, nitro, carboxy, protected carboxy, carbamoyl, carbamoyloxy, cyano, methylsulfonylamino or C<sub>1</sub>-C<sub>4</sub> alkoxy groups. The substituted alkyl groups may be substituted once, twice or three times with the same or with different substituents.

Examples of the above substituted alkyl groups include but are not limited to; cyanomethyl, nitromethyl, hydroxymethyl, trityloxymethyl, propionyloxymethyl, aminomethyl, carboxymethyl, alkyloxycarbonylmethyl, allyloxycarbonylaminomethyl, carbamoyloxymethyl, methoxymethyl, ethoxymethyl, t-butoxymethyl, acetoxymethyl, chloromethyl, bromomethyl, iodomethyl, trifluoromethyl, 6-hydroxyhexyl, 2,4-dichloro(n-butyl), 2-amino(iso-propyl), 2-carbamoyloxyethyl and the like. A preferred group of examples within the above "C<sub>1</sub>-C<sub>12</sub> substituted alkyl" group includes the substituted methyl group, e.g. a methyl group substituted by the same substituents as the "substituted C<sub>n</sub>-C<sub>m</sub> alkyl" group. Examples of the substituted methyl group include groups such as hydroxymethyl, protected hydroxymethyl (e.g. tetrahydropyranyloxymethyl), acetoxymethyl, carbamoyloxymethyl, trifluoromethyl, chloromethyl, bromomethyl and iodomethyl.

The terms "C<sub>1</sub>-C<sub>12</sub> alkyloxy" or "C<sub>1</sub>-C<sub>12</sub> alkoxy" are used interchangeably herein and denote groups such as methoxy, ethoxy, n-propoxy, isopropoxy, n-butoxy, t-butoxy and like groups.

The terms "C<sub>1</sub>-C<sub>12</sub> acyloxy" or "C<sub>1</sub>-C<sub>12</sub> alkanoyloxy" are used interchangeably and denote herein groups such as formyloxy, acetoxy, propionyloxy, butyryloxy, pentanoyloxy, hexanoyloxy, heptanoyloxy, and the like.

The terms "C<sub>1</sub>-C<sub>12</sub> alkylcarbonyl", "C<sub>1</sub>-C<sub>12</sub> alkanoyl" and "C<sub>1</sub>-C<sub>12</sub> acyl" are used interchangeably herein encompass groups such as formyl, acetyl, propionyl, butyryl, pentanoyl, hexanoyl, heptanoyl, benzoyl and the like.

The term "cycloalkyl" as used herein refers to a mono-, bi-, or tricyclic aliphatic ring having 3 to 14 carbon atoms and preferably 3 to 7 carbon atoms.

The term "alkenyl" means a branched or unbranched hydrocarbon radical having the number of carbon atoms designated containing one or more carbon-carbon double bonds, each double bond being independently cis, trans, or a nongeometric isomer.

The term "alkynyl" means a branched or unbranched hydrocarbon radical having the number of carbon atoms designated containing one or more carbon-carbon triple bonds.

The terms "C<sub>1</sub>-C<sub>12</sub> alkylthio" and "C<sub>1</sub>-C<sub>12</sub> substituted alkylthio" denote C<sub>1</sub>-C<sub>12</sub> alkyl and C<sub>1</sub>-C<sub>12</sub> substituted alkyl groups, respectively, attached to a sulfur which is in turn the point of attachment for the alkylthio or substituted alkylthio group to the group or substituent designated.

The term "aryl" when used alone means a homocyclic aromatic radical whether or not fused having the number of carbon atoms designated. Preferred aryl groups include phenyl, naphthyl, biphenyl, phenanthrenyl, naphthacenyl, and the like (see e.g. *Lang's Handbook of Chemistry* (Dean, J. A., ed) 13<sup>th</sup> ed. Table 7-2 [1985]).

The term "substituted phenyl" or "substituted aryl" denotes a phenyl group or aryl group substituted with one, two or three substituents chosen from halogen(F, Cl, Br, I), hydroxy, protected hydroxy, cyano, nitro, C<sub>1</sub>-C<sub>6</sub> alkyl, C<sub>1</sub>-C<sub>6</sub> alkoxy, carboxy, protected carboxy, carboxymethyl, protected carboxymethyl, hydroxymethyl, protected hydroxymethyl, aminomethyl, protected aminomethyl, trifluoromethyl N-(methylsulfonylamino) or other groups specified.

Examples of the term "substituted phenyl" includes but is not limited to a mono- or di(halo)phenyl group such as 4-chlorophenyl, 2,6-dichlorophenyl, 2,5-dichlorophenyl, 3,4-dichlorophenyl, 3-chlorophenyl, 3-bromophenyl, 4-bromophenyl, 3,4-dibromophenyl, 3-chloro-4-fluorophenyl, 2-fluorophenyl and the like; a mono- or di(hydroxy)phenyl group such as 4-hydroxyphenyl, 3-hydroxyphenyl, 2,4-dihydroxyphenyl, the protected-hydroxy derivatives thereof and the like; a nitrophenyl group such as 3- or 4-nitrophenyl; a cyanophenyl group, for example, 4-cyanophenyl; a mono- or di(lower alkyl)phenyl group such as 4-methylphenyl, 2,4-dimethylphenyl, 2-methylphenyl, 4-(iso-propyl)phenyl, 4-ethylphenyl, 3-(n-propyl)phenyl and the like; a mono or di(alkoxy)phenyl group, for example, 2,6-dimethoxyphenyl, 4-methoxyphenyl, 3-ethoxyphenyl, 4-(isopropoxy)phenyl, 4-(t-butoxy)phenyl, 3-ethoxy-4-methoxyphenyl and the like; 3- or 4- trifluoromethylphenyl; a mono- or dicarboxyphenyl or (protected carboxy)phenyl group such 4-carboxyphenyl; a mono- or di(hydroxymethyl)phenyl or (protected hydroxymethyl)phenyl such as 3-(protected hydroxymethyl)phenyl or 3,4-di(hydroxymethyl)phenyl; a mono- or di(aminomethyl)phenyl or (protected aminomethyl)phenyl such as 2-(aminomethyl)phenyl or 2,4-(protected aminomethyl)phenyl; or a mono- or di(N-(methylsulfonylamino))phenyl such as 3-(N-methylsulfonylamino))phenyl. Also, the term "substituted phenyl" represents disubstituted phenyl groups wherein the substituents are different, for example, 3-methyl-4-hydroxyphenyl, 3-chloro-4-hydroxyphenyl, 2-methoxy-4-bromophenyl, 4-ethyl-2-hydroxyphenyl, 3-hydroxy-4-nitrophenyl, 2-hydroxy-4-chlorophenyl and the like. Preferred substituted phenyl groups include the 2- and 3-trifluoromethylphenyl, the 4-hydroxyphenyl, the 2-aminomethylphenyl and the 3-(N-(methylsulfonylamino))phenyl groups.



The term "arylalkyl" means one, two, or three aryl groups having the number of carbon atoms designated, appended to an alkyl radical having the number of carbon atoms designated including but not limited to; benzyl, naphthylmethyl, phenethyl, benzhydryl (diphenylmethyl), trityl, and the like. A preferred arylalkyl group is the benzyl group.

5       The term "substituted C<sub>6</sub>-C<sub>10</sub>aryl-C<sub>1</sub>-C<sub>8</sub>alkyl" denotes a C<sub>1</sub>-C<sub>8</sub>alkyl group substituted at any carbon with a C<sub>6</sub>-C<sub>10</sub>aryl group bonded to the alkyl group through any aryl ring position and substituted on the C<sub>1</sub>-C<sub>8</sub>alkyl portion with one, two or three groups chosen from halogen (F, Cl, Br, I), hydroxy, protected hydroxy, amino, protected amino, C<sub>1</sub>-C<sub>7</sub>acyloxy, nitro, carboxy, protected carboxy, carbamoyl, carbamoyloxy, cyano, C<sub>1</sub>-C<sub>6</sub>alkylthio, N-  
10 (methylsulfonylamino) or C<sub>1</sub>-C<sub>4</sub>alkoxy. Optionally the aryl group may be substituted with one, two, or three groups chosen from halogen, hydroxy, protected hydroxy, nitro, C<sub>1</sub>-C<sub>6</sub>alkyl, C<sub>1</sub>-C<sub>4</sub>alkoxy, carboxy, protected carboxy, carboxymethyl, protected carboxymethyl, hydroxymethyl, protected hydroxymethyl, aminomethyl, protected aminomethyl, or an N-(methylsulfonylamino) group. As before, when either the C<sub>1</sub>-C<sub>8</sub>alkyl portion or the aryl  
15 portion or both are disubstituted, the substituents can be the same or different.

Examples of the term "substituted C<sub>6</sub>-C<sub>10</sub>aryl-C<sub>1</sub>-C<sub>8</sub>alkyl" include groups such as 2-phenyl-1-chloroethyl, 2-(4-methoxyphenyl)ethyl, 2,6-dihydroxy-4-phenyl(n-hexyl), 5-cyano-3-methoxy-2-phenyl(n-pentyl), 3-(2,6-dimethylphenyl)n-propyl, 4-chloro-3-aminobenzyl, 6-(4-methoxyphenyl)-3-carboxy(n-hexyl), 5-(4-aminomethyl phenyl)-3-(aminomethyl)(n-pentyl),  
20 and the like.

The term "carboxy-protecting group" as used herein refers to one of the ester derivatives of the carboxylic acid group commonly employed to block or protect the carboxylic acid group while reactions are carried out on other functional groups on the compound. Examples of such carboxylic acid protecting groups include 4-nitrobenzyl, 4-methoxybenzyl,  
25 3,4-dimethoxybenzyl, 2,4-dimethoxybenzyl, 2,4,6-trimethoxybenzyl, 2,4,6-trimethylbenzyl, pentamethylbenzyl, 3,4-methylenedioxybenzyl, benzhydryl, 4,4'-dimethoxybenzhydryl, 2,2',4,4'-tetramethoxybenzhydryl, t-butyl, t-amyl, trityl, 4-methoxytrityl, 4,4'-dimethoxytrityl, 4,4',4''-trimethoxytrityl, 2-phenylprop-2-yl, trimethylsilyl, t-butyl dimethylsilyl, phenacyl, 2,2,2-trichloroethyl,  $\beta$ -(trimethylsilyl)ethyl,  $\beta$ -(di(n-butyl)methylsilyl)ethyl, p-toluenesulfonyl ethyl, 4-nitrobenzylsulfonyl ethyl, allyl, cinnamyl, 1-(trimethylsilylmethyl)prop-1-en-3-yl, and like  
30 moieties. The species of carboxy-protecting group employed is not critical so long as the derivatized carboxylic acid is stable to the condition of subsequent reaction(s) on other positions of the benzodiazepinedione molecule and can be removed at the appropriate point without disrupting the remainder of the molecule. In particular, it is important not to subject  
35 the carboxy-protected benzodiazepinedione molecule to strong nucleophilic bases or reductive conditions employing highly activated metal catalysts such as Raney nickel. (Such harsh removal conditions are also to be avoided when removing amino-protecting groups and hydroxy-protecting groups, discussed below.) Preferred carboxylic acid protecting groups are the allyl and p-nitrobenzyl groups. Similar carboxy-protecting groups used in the

cephalosporin, penicillin and peptide arts can also be used to protect a carboxy group substituents of the benzodiazepinedione. Further examples of these groups are found in E. Haslam, "Protective Groups in Organic Chemistry", J. G. W. McOmie, Ed., Plenum Press, New York, N.Y., 1973, Chapter 5, and T.W. Greene, "Protective Groups in Organic Synthesis", John Wiley and Sons, New York, NY, 1981, Chapter 5. The term "protected carboxy" refers to a carboxy group substituted with one of the above carboxy-protecting groups.

As used herein the term "amide-protecting group" refers to any group typically used in the peptide art for protecting the peptide nitrogens from undesirable side reactions. Such groups include p-methoxyphenyl, 3,4-dimethoxybenzyl, benzyl, O-nitrobenzyl, di-(p-methoxyphenyl)methyl, triphenylmethyl, (p-methoxyphenyl)diphenylmethyl, diphenyl-4-pyridylmethyl, m-2-(picolyl)-N'-oxide, 5-dibenzosuberyl, trimethylsilyl, t-butyl dimethylsilyl, and the like. Further descriptions of these protecting groups can be found in "Protective Groups in Organic Synthesis", by Theodora W. Greene, 1981, John Wiley and Sons, New York.

Unless otherwise specified, the terms "heterocyclic group" or "heterocyclic" or "HET" or "heterocyclyl" are used interchangeably as used herein refer to any mono-, bi-, or tricyclic saturated, unsaturated, or aromatic ring having the number of atoms designated where at least one ring is a 5-, 6- or 7-membered ring containing from one to four heteroatoms selected from the group nitrogen, oxygen, and sulfur preferably at least one heteroatom is nitrogen (*Lang's Handbook of Chemistry, supra*). Typically, the 5-membered ring has 0 to 2 double bonds and the 6- or 7-membered ring has 0 to 3 double bonds and the nitrogen or sulfur heteroatoms may optionally be oxidized, and any nitrogen heteroatom may optionally be quaternized. Included in the definition are any bicyclic groups where any of the above heterocyclic rings are fused to a benzene ring. Heterocyclics in which nitrogen is the heteroatom are preferred.

The following ring systems are examples of the heterocyclic (whether substituted or unsubstituted) radicals denoted by the term "heterocyclic": thienyl, furyl, pyrrolyl, imidazolyl, pyrazolyl, thiazolyl, isothiazolyl, oxazolyl, isoxazolyl, triazolyl, thiadiazolyl, oxadiazolyl, tetrazolyl, thiatriazolyl, oxatriazolyl, pyridyl, pyrimidyl, pyrazinyl, pyridazinyl, thiazinyl, oxazinyl, triazinyl, thiadiazinyl, oxadiazinyl, dithiazinyl, dioxazinyl, oxathiazinyl, tetrazinyl, thiatriazinyl, oxatriazinyl, dithiadiazinyl, imidazoliny, dihydropyrimidyl, tetrahydropyrimidyl, tetrazolo[1,5-b]pyridazinyl and purinyl, as well as benzo-fused derivatives, for example benzoxazolyl, benzthiazolyl, benzimidazolyl and indolyl.

Heterocyclic 5-membered ring systems containing a sulfur or oxygen atom and one to three nitrogen atoms are also suitable for use in the instant invention. Examples of such preferred groups include thiazolyl, in particular thiazol-2-yl and thiazol-2-yl N-oxide, thiadiazolyl, in particular 1,3,4-thiadiazol-5-yl and 1,2,4-thiadiazol-5-yl, oxazolyl, preferably oxazol-2-yl, and oxadiazolyl, such as 1,3,4-oxadiazol-5-yl, and 1,2,4-oxadiazol-5-yl. A group of further preferred examples of 5-membered ring systems with 2 to 4 nitrogen atoms include imidazolyl, preferably imidazol-2-yl; triazolyl, preferably 1,3,4-triazol-5-yl; 1,2,3-triazol-5-yl,

1,2,4-triazol-5-yl, and tetrazolyl, preferably 1H-tetrazol-5-yl. A preferred group of examples of benzo-fused derivatives are benzoxazol-2-yl, benzthiazol-2-yl and benzimidazol-2-yl.

Further suitable specific examples of the above heterocyclic ring systems are 6-membered ring systems containing one to three nitrogen atoms. Such examples include pyridyl, such as pyrid-2-yl, pyrid-3-yl, and pyrid-4-yl; pyrimidyl, preferably pyrimid-2-yl and pyrimid-4-yl; triazinyl, preferably 1,3,4-triazin-2-yl and 1,3,5-triazin-4-yl; pyridazinyl, in particular pyridazin-3-yl, and pyrazinyl. The pyridine N-oxides and pyridazine N-oxides and the pyridyl, pyrimid-2-yl, pyrimid-4-yl, pyridazinyl and the 1,3,4-triazin-2-yl radicals, are a preferred group.

The substituents for the optionally substituted heterocyclic ring systems, and further examples of the 5- and 6-membered ring systems discussed above can be found in W. Druckheimer *et al.*, U.S. Patent No. 4,278,793.

A particularly preferred group of "heterocyclics" or "HET" include; 1,3-thiazol-2-yl, 4-(carboxymethyl)-5-methyl-1,3-thiazol-2-yl, 4-(carboxymethyl)-5-methyl-1,3-thiazol-2-yl sodium salt, 1,2,4-thiadiazol-5-yl, 3-methyl-1,2,4-thiadiazol-5-yl, 1,3,4-triazol-5-yl, 2-methyl-1,3,4-triazol-5-yl, 2-hydroxy-1,3,4-triazol-5-yl, 2-carboxy-4-methyl-1,3,4-triazol-5-yl sodium salt, 2-carboxy-4-methyl-1,3,4-triazol-5-yl, 1,3-oxazol-2-yl, 1,3,4-oxadiazol-5-yl, 2-methyl-1,3,4-oxadiazol-5-yl, 2-(hydroxymethyl)-1,3,4-oxadiazol-5-yl, 1,2,4-oxadiazol-5-yl, 1,3,4-thiadiazol-5-yl, 2-thiol-1,3,4-thiadiazol-5-yl, 2-(methylthio)-1,3,4-thiadiazol-5-yl, 2-amino-1,3,4-thiadiazol-5-yl, 1H-tetrazol-5-yl, 1-methyl-1H-tetrazol-5-yl, 1-(1-(dimethylamino)eth-2-yl)-1H-tetrazol-5-yl, 1-(carboxymethyl)-1H-tetrazol-5-yl, 1-(carboxymethyl)-1H-tetrazol-5-yl sodium salt, 1-(methylsulfonic acid)-1H-tetrazol-5-yl, 1-(methylsulfonic acid)-1H-tetrazol-5-yl sodium salt, 2-methyl-1H-tetrazol-5-yl, 1,2,3-triazol-5-yl, 1-methyl-1,2,3-triazol-5-yl, 2-methyl-1,2,3-triazol-5-yl, 4-methyl-1,2,3-triazol-5-yl, pyrid-2-yl N-oxide, 6-methoxy-2-(n-oxide)-pyridaz-3-yl, 6-hydroxypyridaz-3-yl, 1-methylpyrid-2-yl, 1-methylpyrid-4-yl, 2-hydroxypyrimid-4-yl, 1,4,5,6-tetrahydro-5,6-dioxo-4-methyl-as-triazin-3-yl, 1,4,5,6-tetrahydro-4-(formylmethyl)-5,6-dioxo-as-triazin-3-yl, 2,5-dihydro-5-oxo-6-hydroxy-as-triazin-3-yl, 2,5-dihydro-5-oxo-6-hydroxy-as-triazin-3-yl sodium salt, 2,5-dihydro-5-oxo-6-hydroxy-2-methyl-as-triazin-3-yl sodium salt, 2,5-dihydro-5-oxo-6-hydroxy-2-methyl-as-triazin-3-yl, 2,5-dihydro-5-oxo-6-methoxy-2-methyl-as-triazin-3-yl, 2,5-dihydro-5-oxo-as-triazin-3-yl, 2,5-dihydro-5-oxo-2-methyl-as-triazin-3-yl, 2,5-dihydro-5-oxo-2,6-dimethyl-as-triazin-3-yl, tetrazolo[1,5-b]pyridazin-6-yl and 8-aminotetrazolo[1,5-b]-pyridazin-6-yl.

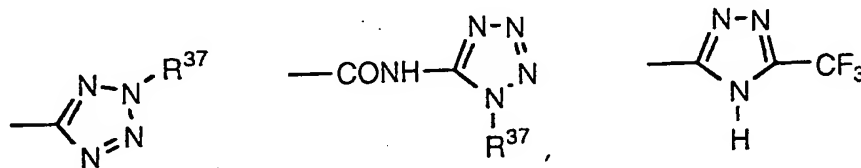
An alternative group of "heterocyclics" includes; 4-(carboxymethyl)-5-methyl-1,3-thiazol-2-yl, 4-(carboxymethyl)-5-methyl-1,3-thiazol-2-yl sodium salt, 1,3,4-triazol-5-yl, 2-methyl-1,3,4-triazol-5-yl, 1H-tetrazol-5-yl, 1-methyl-1H-tetrazol-5-yl, 1-(1-(dimethylamino)eth-2-yl)-1H-tetrazol-5-yl, 1-(carboxymethyl)-1H-tetrazol-5-yl, 1-(carboxymethyl)-1H-tetrazol-5-yl sodium salt, 1-(methylsulfonic acid)-1H-tetrazol-5-yl, 1-(methylsulfonic acid)-1H-tetrazol-5-yl sodium salt, 1,2,3-triazol-5-yl, 1,4,5,6-tetrahydro-5,6-dioxo-4-methyl-as-triazin-3-yl, 1,4,5,6-tetrahydro-4-(2-formylmethyl)-5,6-dioxo-as-triazin-3-yl, 2,5-dihydro-5-oxo-6-hydroxy-2-

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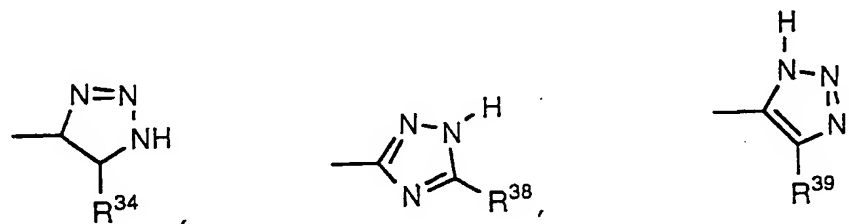
methyl-as-triazin-3-yl sodium salt, 2,5-dihydro-5-oxo-6-hydroxy-2-methyl-as-triazin-3-yl, tetrazolo[1,5-b]pyridazin-6-yl, and 8-aminotetrazolo[1,5-b]pyridazin-6-yl.

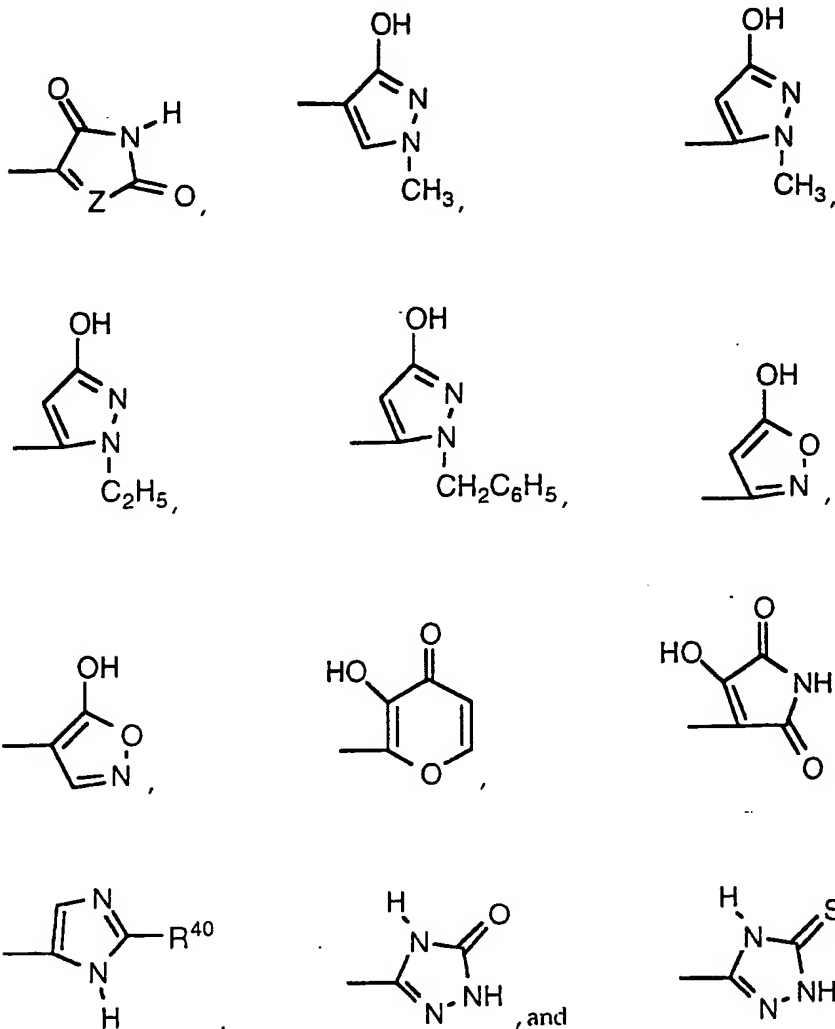
Bivalent radicals L (either L<sup>1</sup>, L<sup>2</sup>, or L<sup>3</sup>) whether branched or unbranched, derived from alkanes, alkenes, alkadienes, alkynes, alkadiynes, and arenes optionally containing O, N and/or S atoms, or homo- and heterocycles either aromatic or aliphatic, are designated by adding the suffix "ene" to the corresponding monovalent radical. Atoms bearing the free valences may include any C, O, N or S.

The phrase "negatively charged acidic moiety" as used to define the W moiety, is intended to embrace chemical groups which, when attached to the 7-member lactam ring of Formula I, confers acidic character to the compound of Formula I. "Acidic character" means proton-donor capability, that is, the capacity of the compound of Formula I to be a proton donor in the presence of a proton acceptor, such as water. Typically, the acidic group should be selected to have proton donor capability such that the compound of Formula I has a pK<sub>a</sub> in a range from about one to about twelve. More typically, the Formula I compound would have a pK<sub>a</sub> in a range from about two to about seven. Also typically, W will be a group that contains at least one acidic hydrogen. Optionally, W will be a prodrug that is capable of being converted, *in vivo*, into a negatively charged acidic moiety. An example of an acidic group containing at least one acidic hydrogen atom is the carboxyl group (-COOH). A preferred example of W in Formula I containing a carboxyl group is where R<sup>27</sup> is -CH<sub>2</sub>- and "w" is -COOH. There are many examples of equivalent acidic groups other than carboxyl group, that can be substituted into W to produce biologically equivalent compounds of Formula I. Such other acidic groups are collectively referred to as "bioisosteres of the carboxylate" or "acidic bioisosteres". Representative examples of such "basic bioisosteres" are set forth above under "w" (a) through (ap). Other specific examples of such acidic bioisosteres may be independently selected from acidic moieties consisting of CO<sub>2</sub>H, CO<sub>2</sub>CH<sub>3</sub>, SH, CH<sub>2</sub>SH, C<sub>2</sub>H<sub>4</sub>SH, PO<sub>3</sub>H<sub>2</sub>, NHSO<sub>2</sub>CF<sub>3</sub>, NHSO<sub>2</sub>C<sub>6</sub>F<sub>5</sub>, SO<sub>3</sub>H, CONHNH<sub>2</sub>, CONHNHSO<sub>2</sub>CF<sub>3</sub>, CONHOCH<sub>3</sub>, CONHOC<sub>2</sub>H<sub>5</sub>, CONHCF<sub>3</sub>, OH, CH<sub>2</sub>OH, C<sub>2</sub>H<sub>4</sub>OH, OPO<sub>3</sub>H<sub>2</sub>, OSO<sub>3</sub>H,



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wherein each of  $R^{37}$ ,  $R^{38}$ ,  $R^{39}$ , and  $R^{40}$  is independently selected from H, Cl, CN,  $NO_2$ ,  $CF_3$ ,  $C_2F_5$ ,  $C_3F_7$ ,  $CHF_2$ ,  $CH_2F$ ,  $CO_2CH_3$ ,  $CO_2C_2H_5$ ,  $SO_2CH_3$ ,  $SO_2CF_3$  and  $SO_2C_6H_5$ ; wherein Z is selected from O, S,  $NR^{41}$  and  $CH_2$ ; wherein  $R^{41}$  is selected from hydrido,  $CH_3$  and  $CH_2C_6H_5$ ; or a tautomer thereof or a pharmaceutically-acceptable salt thereof.

"Pharmaceutically acceptable salts" include both acid and base addition salts.

"Pharmaceutically acceptable acid addition salt" refers to those salts which retain the biological effectiveness and properties of the free bases and which are not biologically or otherwise undesirable, formed with inorganic acids such as hydrochloric acid, hydrobromic acid, sulfuric acid, nitric acid, carbonic acid, phosphoric acid and the like, and organic acids may be selected from aliphatic, cycloaliphatic, aromatic, araliphatic, heterocyclic, carboxylic, and sulfonic classes of organic acids such as formic acid, acetic acid, propionic acid, glycolic acid, gluconic acid, lactic acid, pyruvic acid, oxalic acid, malic acid, maleic acid, malonic acid,

succinic acid, fumaric acid, tartaric acid, citric acid, aspartic acid, ascorbic acid, glutamic acid, anthranilic acid, benzoic acid, cinnamic acid, mandelic acid, embonic acid, phenylacetic acid, methanesulfonic acid, ethanesulfonic acid, p-toluenesulfonic acid, salicylic acid and the like.

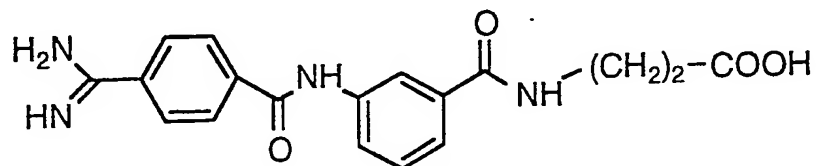
"Pharmaceutically acceptable base addition salts" include those derived from inorganic bases such as sodium, potassium, lithium, ammonium, calcium, magnesium, iron, zinc, copper, manganese, aluminum salts and the like. Particularly preferred are the ammonium, potassium, sodium, calcium and magnesium salts. Salts derived from pharmaceutically acceptable organic nontoxic bases includes salts of primary, secondary, and tertiary amines, substituted amines including naturally occurring substituted amines, cyclic amines and basic ion exchange resins, such as isopropylamine, trimethylamine, diethylamine, triethylamine, tripropylamine, ethanolamine, 2-diethylaminoethanol, trimethylamine, dicyclohexylamine, lysine, arginine, histidine, caffeine, procaine, hydrabamine, choline, betaine, ethylenediamine, glucosamine, methylglucamine, theobromine, purines, piperazine, piperidine, N-ethylpiperidine, polyamine resins and the like. Particularly preferred organic non-toxic bases are isopropylamine, diethylamine, ethanolamine, trimethylamine, dicyclohexylamine, choline, and caffeine.

The term "prodrug" as used herein means a derivative of a parent drug molecule that enhances pharmaceutically desirable characteristics or properties (e.g. transport, bioavailability, pharmacodynamics, etc.) and that requires biotransformation, either spontaneous or enzymatic, within the organism to release the active parent drug.

#### 20 B. Utility

The present invention is the result of the unexpected discovery that substituted fused 7-membered "lactam" ring compounds defined by formula I, especially including substituted 3,4-dihydro-1H-1,4-benzodiazepine-2,5-diones (referred to hereafter as substituted benzodiazepinediones) are capable of acting as fibrinogen receptor antagonists, that is inhibiting both platelet aggregation and binding of fibrinogen to the platelet receptor, GPIIb/IIIa, at concentrations comparable to many of the more potent anti-aggregatory peptides of the prior art (see e.g. Ali *et al.* EP 341 915). Thus, the instant fibrinogen receptor antagonists are particularly useful when they inhibit the binding of fibrinogen to the GPIIb/IIIa receptor with an IC<sub>50</sub> in a Fg/GPIIb/IIIa ELISA as described herein of at least about 35nM and an IC<sub>50</sub> in a platelet aggregation (PA) assay, also described herein, of at least about 3μM. Since there is some variability in the values obtained for ELISA and PA assays in different laboratories, preferred fibrinogen receptor antagonists represented by Formulae I-VI will have IC<sub>50</sub>'s for Fg/GPIIb/IIIa ELISA's and PA assays equal to or greater than that of the prior art compound N-[3-(4-amidinobenzamido)benzoyl]-3-aminopropionic acid,

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This compound is used as an internal standard and has been found to have an Fg/GPIIb/IIIa ELISA IC<sub>50</sub> of about 25 nM and a PA inhibition assay IC<sub>50</sub> of about 3 μM when measured according to the assays described herein. Most preferred fibrinogen receptor antagonists represented by Formulae I-VI have IC<sub>50</sub>s in Fg/GPIIb/IIIa ELISA of at least about 15 nM and in platelet aggregation inhibition assays of at least about 1 μM or three times more potent in a platelet aggregation inhibition assay than the reference compound N-[3-(4-amidinobenzamido)benzoyl]-3-aminopropionic acid. This reference compound is readily prepared by standard methods from commercially available precursors (e.g. β-alanine, *m*-aminobenzoic acid, and *p*-aminobenzoic acid) or may be synthesized according to the procedures set forth in EP 372 486 A2.

The compounds described in the present invention may therefore inhibit the binding of fibrinogen to its receptor on platelets, GPIIb/IIIa, and thus prevent the aggregation of platelets and the formation of platelet plugs, emboli and thrombi in the circulatory system in mammals. Thromboembolic disorders have been shown to be directly related to the susceptibility of blood platelets to aggregate. Mammals exposed to medical procedures such as angioplasty and thrombolytic therapy are particularly susceptible to thrombus formation. The compounds of the present invention can be used to inhibit thrombus formation following angioplasty. They may also be used in combination with antithrombolytic agents such as tissue plasminogen activator and its derivatives (US patents 4,752,603; 4,766,075; 4,777,043; EP 199 574; EP 238 304; EP 228 862; EP 297 860; PCT WO89/04368; PCT WO89/00197), streptokinase and its derivatives, or urokinase and its derivatives to prevent arterial reocclusion following thrombolytic therapy. When used in combination with the above thrombolytic agents, the compounds of the present invention may be administered prior to, simultaneously with, or subsequent to the antithrombolytic agent. Mammals exposed to renal dialysis, blood oxygenation, cardiac catheterization and similar medical procedures as well as mammals fitted with certain prosthetic devices are also susceptible to thromboembolic disorders. Physiologic conditions, with or without known cause may also lead to thromboembolic disorders. Thus, the compounds described herein may be useful in treating thromboembolic disorders in mammals. The compounds described herein may also be used as adjuncts to anticoagulant therapy, for example in combination with aspirin, heparin or warfarin and other anticoagulant agents. The application of the compounds described herein for these and related disorders will be apparent to those skilled in the art.

Compounds represented by Formula I are also capable of antagonizing the interaction between vitronectin(Vn) and the vitronectin receptor (VnR) and thus are useful in situations where such antagonism is indicated. Vn-VnR antagonism is particularly pronounced when substituent Q<sup>1</sup>-L<sup>1</sup>- is bonded to position 8 or 9 of the substituted benzodiazapinedione, as represented by Formula VIa below. Thus, it is contemplated the instant compounds are useful in treating bone disorders such as osteoporosis.

Compounds of this invention are also useful as intermediates generally, or as precursors of inhibitors to GPIIb/IIIa or other integrin receptors and thus in addition to treating cardiovascular and bone disease, these compounds may be usefully employed in metastatic disease, or for any disease where antagonism of RGD binding integrin receptors is indicated.

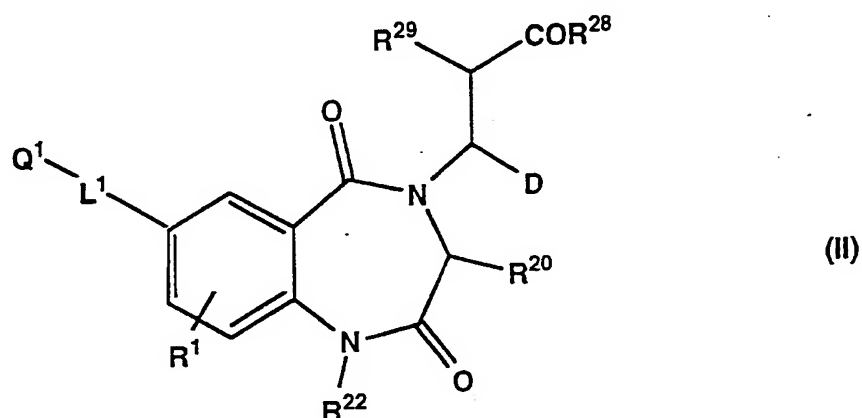
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### C. Preferred Embodiments

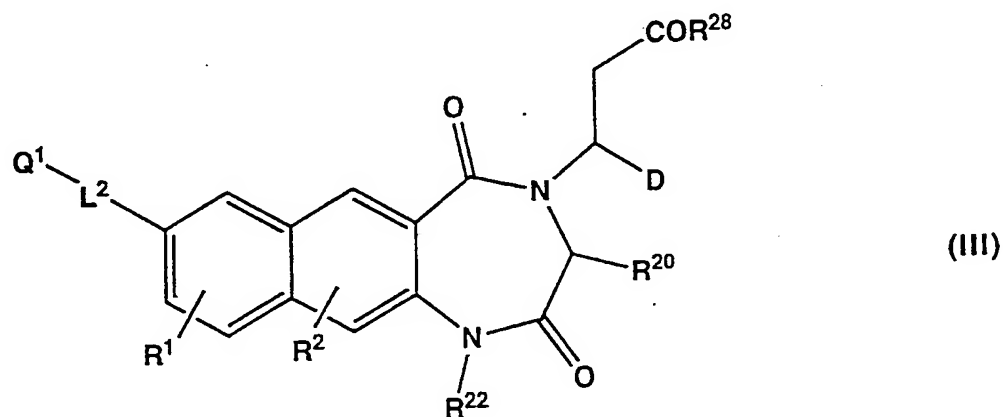
#### 1. Nonpeptidyl GPIIb/IIIa Inhibitors

One embodiment of the invention comprises a compound represented by formula I capable of inhibiting binding of the platelet GPIIb/IIIa receptor to its native *in vivo* ligands.

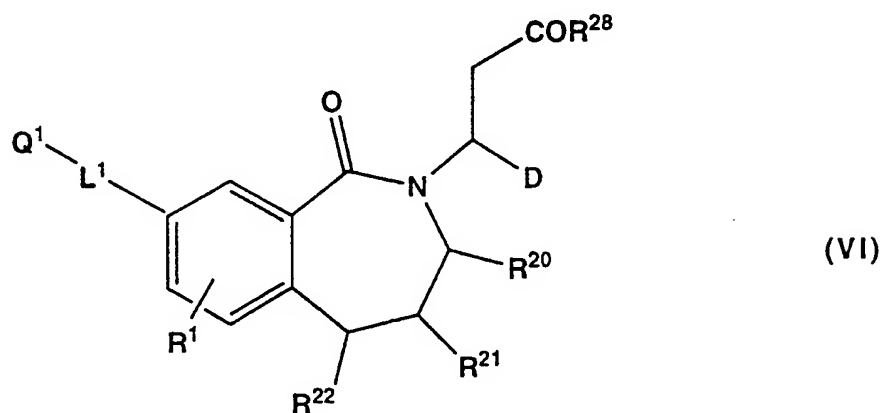
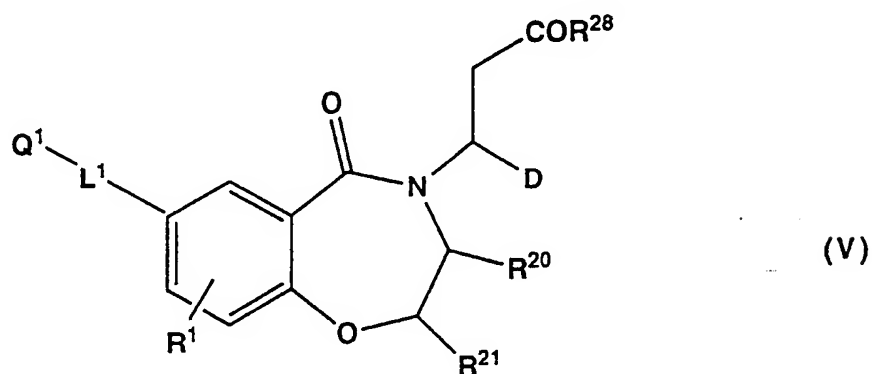
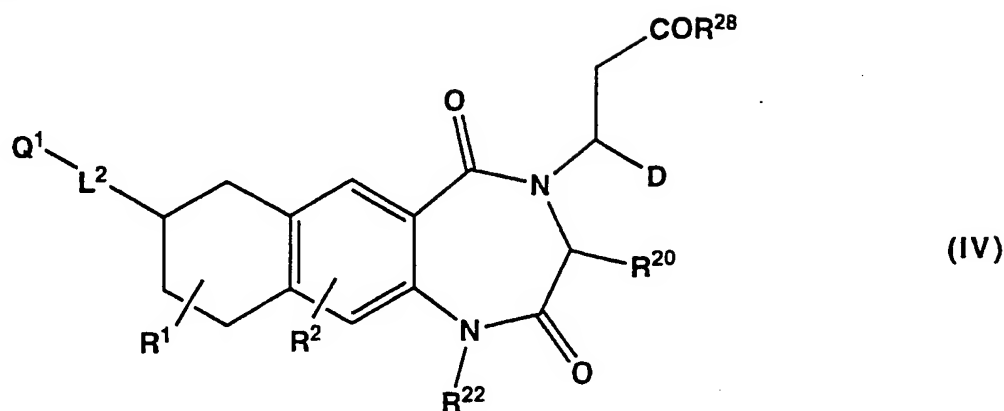
10 Preferred nonpeptidyl inhibitors include compounds represented by structural formulae II-VI:



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where  $R^1, R^2, R^{20}, R^{21}, R^{22}, R^{28}, R^{29}, Q^1, L^1$  and  $L^2$  are defined above.

Referring to formulae I-VI the following structural features of the instant nonpeptidyl antithrombotic inhibitors can be identified:

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- The positively charged Q moiety;
- The linking moiety L;
- The flat (usually aromatic) ring A;
- The 7-member "lactam" ring containing TUG;
- Substituents of TUG especially  $R^{22}$ ;

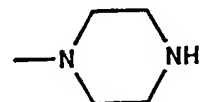
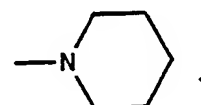
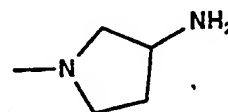
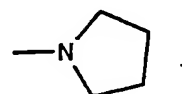
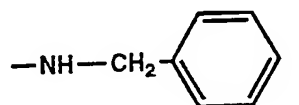
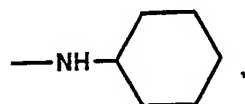
- f. The amino acid linking moiety D; and  
g. The negatively charged acidic moiety W.

a. Positively charged Q

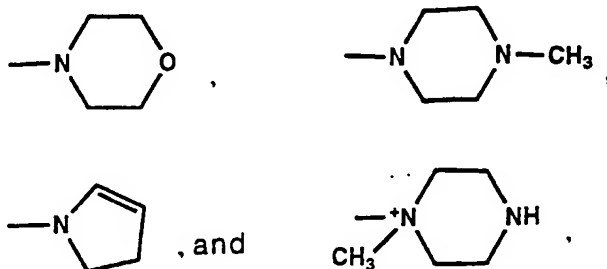
Suitable groups Q (either Q<sup>1</sup> or Q<sup>2</sup>) contain one or more nitrogen atoms and have a pK<sub>b</sub> sufficiently high so that they are at least 10% positively charged at physiological pH. Q may be one or more primary, secondary, tertiary, or quaternary amines or imines either isolated or conjugated with other nitrogen atoms to form groups including but not limited to; aminomethyleneimino, amidino, and guanidino groups and multiples thereof. Alternatively, Q may be a saturated or unsaturated (including aromatic) heterocyclic group provided the group bears a positive charge at physiological pH. In one embodiment of the invention, Q is preferably selected from; amino (H<sub>2</sub>N-), imino (=NH), amidino (H<sub>2</sub>NC(=NH)-), aminomethyleneamino (H<sub>2</sub>N-CH=N-), iminomethylamino (HN=CH-NH-), guanidino (H<sub>2</sub>N-C(=NH)-NH-), N<sup>G</sup>-aminoguanidino (H<sub>2</sub>N-HN-C(=NH)-NH-), alkylamino (R<sup>1</sup>NH-), dialkylamino (R<sup>1</sup><sub>2</sub>N-), trialkylamino (R<sup>1</sup><sub>3</sub>N-), alkylideneamino (R<sup>1</sup><sub>2</sub>C=N-), pyranyl, pyrrolyl, imadazolyl, pyrazolyl, pyridyl, pyrazinyl, pyrimidinyl, pyridazinyl, indoliziny, isoindolyl, 3H-indolyl, indolyl, 1H-indazolyl, purinyl, 4H-quinoliziny, isoquinolyl, quinolyl, phthalazinyl, naphthyridinyl, quinoxaliny, quinazolinyl, cinnoliny, pteridinyl, 4aH-carbazolyl, carbazolyl, b-carboliny, phenanthridinyl, acridinyl, perimidinyl, phenanthroliny, phenazinyl, phenarsazinyl, phenothiazinyl, pyrrolinyl, imidazolidinyl, imidazoliny, pyrazolidinyl, pyrazolinyl, piperidyl, piperazinyl, indoliny, isoindoliny, quinuclidinyl, morpholiny, and 1,3-diazacyclohex-4-ene, where R<sup>1</sup> is selected from; hydrogen, C<sub>1</sub>-C<sub>6</sub>-alkyl, C<sub>6</sub>-C<sub>10</sub>-aryl, and C<sub>3</sub>-C<sub>10</sub>-cycloalkyl. Optionally, any of the nitrogen containing heterocycles described above may be substituted with amino, imino, amidino, aminomethyleneamino, iminomethylamino, guanidino, N<sup>G</sup>-amino-guanidino, alkylamino, dialkylamino, trialkylamino, or alkylideneamino groups.

Exemplary preferred Q<sup>1</sup> groups include the following:

(A) Amino and ammonio groups;

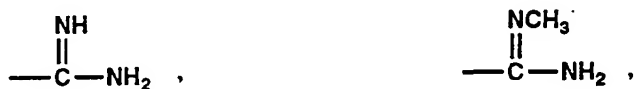


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(B) Amidino groups;

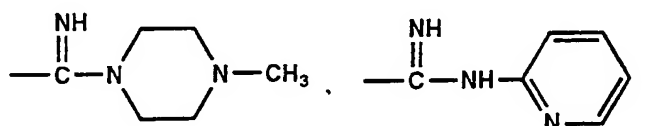
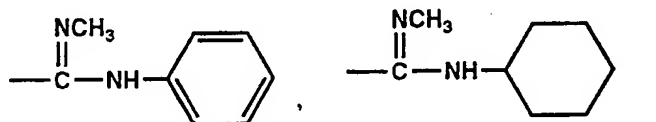
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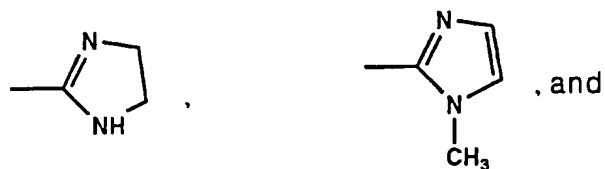
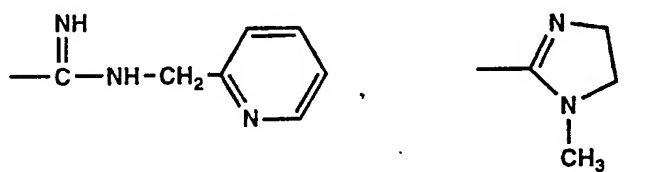
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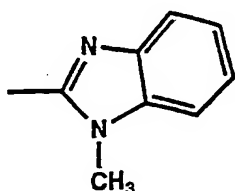


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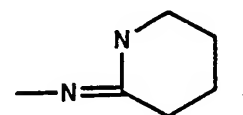
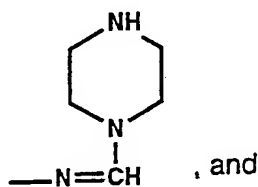
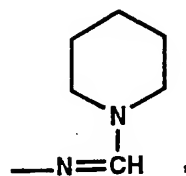
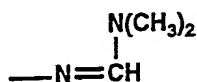
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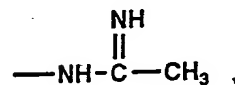
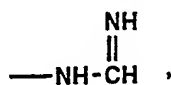
(C) Aminoalkyleneimino groups;

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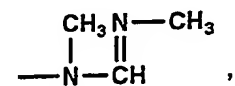
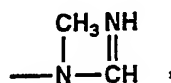


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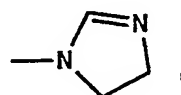
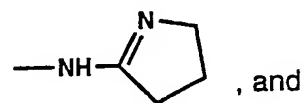
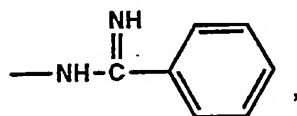
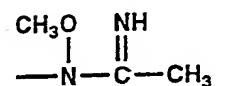
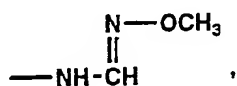
(D) Iminoalkyleneamino groups;



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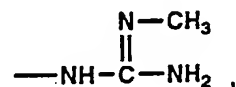
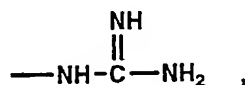


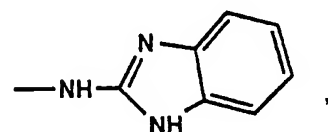
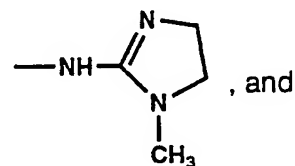
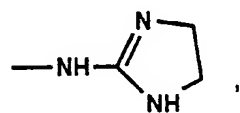
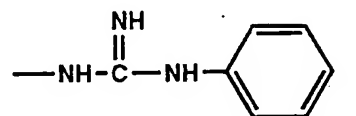
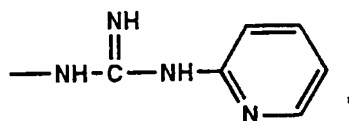
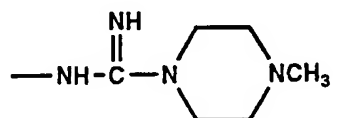
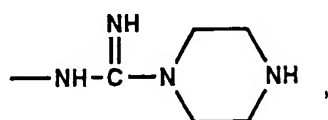
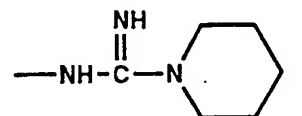
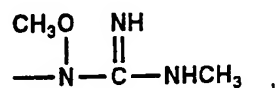
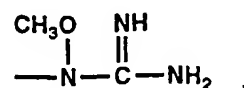
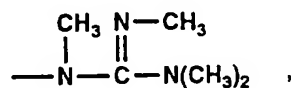
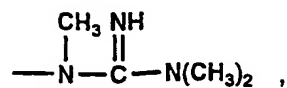
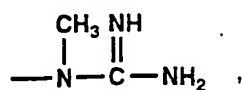
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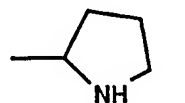
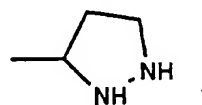
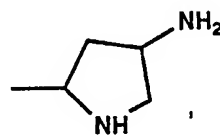
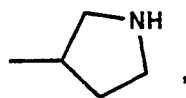
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(E) Guanidino groups;

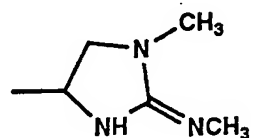
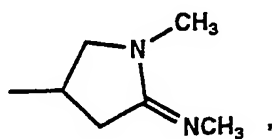
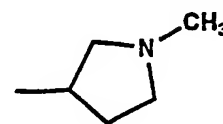
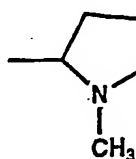




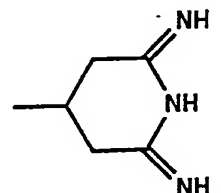
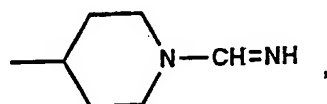
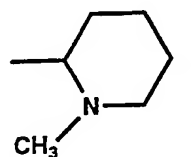
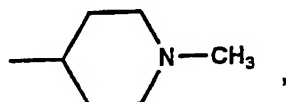
(F) Saturated heterocyclic groups;



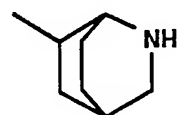
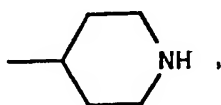
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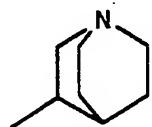
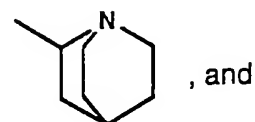
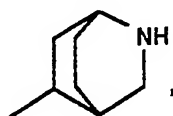
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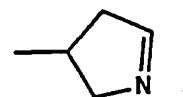
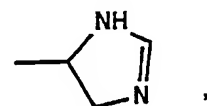


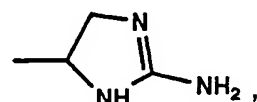
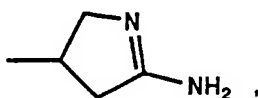
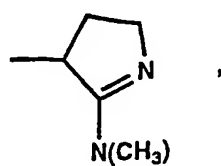
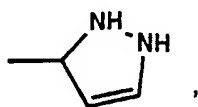
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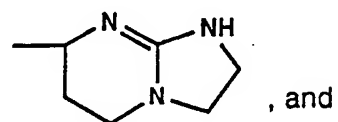
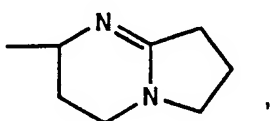
(G) Unsaturated (nonaromatic) heterocyclic groups;

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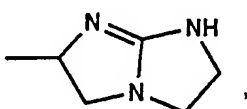




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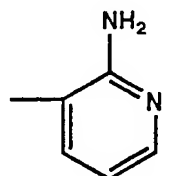
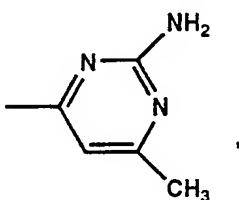


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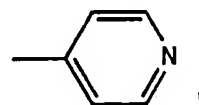
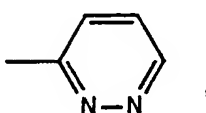
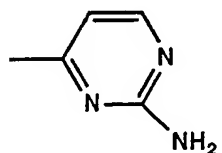
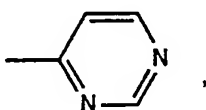


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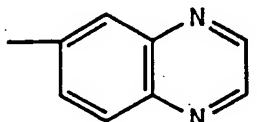
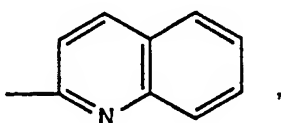
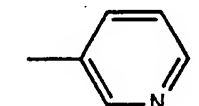
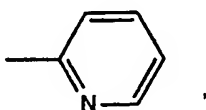
(H) Unsaturated aromatic heterocyclic groups:



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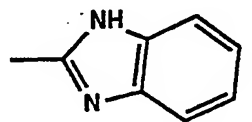
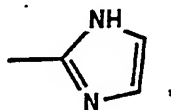
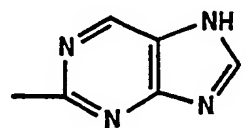
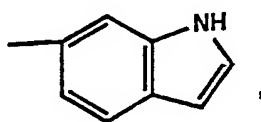


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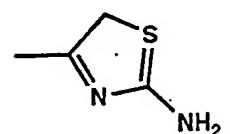
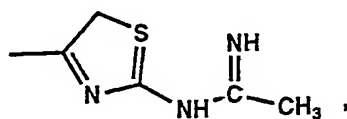
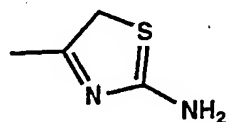
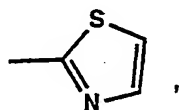


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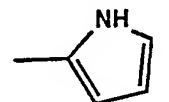
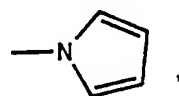
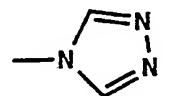
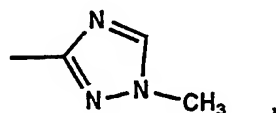
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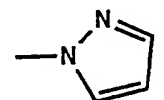
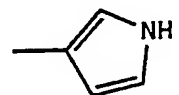
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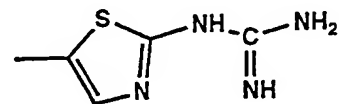
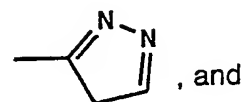
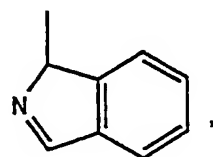
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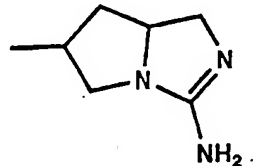
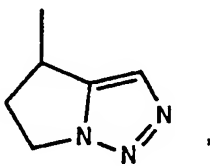
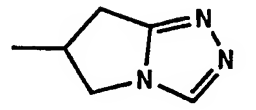
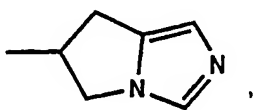
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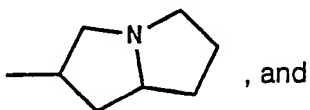
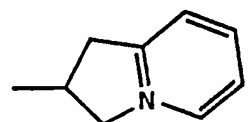
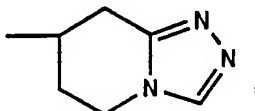


## (I) Bicycloheterocyclic groups;

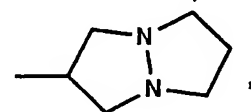
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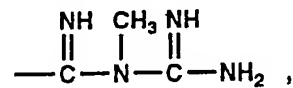
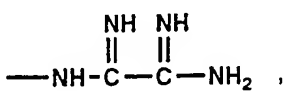


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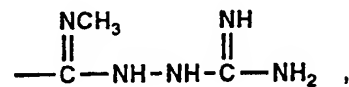
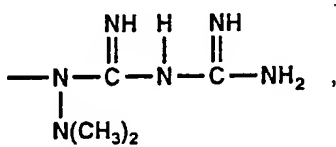
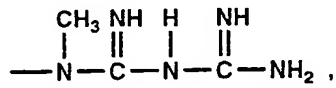
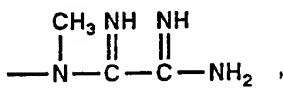


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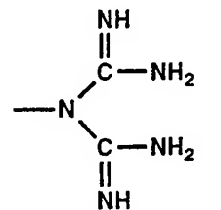
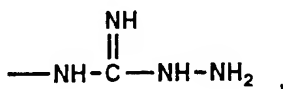
## (J) Multiple amino, guanidino, and amidino groups;

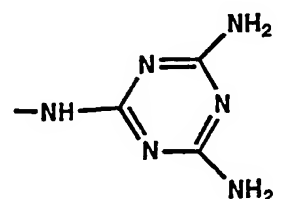
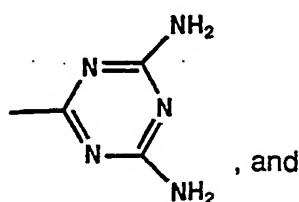


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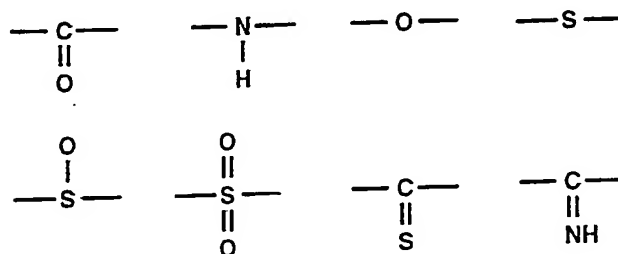




Most preferred Q groups are amino, amidino, and guanidino groups.

b. The linking group L

- 5        The length of the bivalent radical L appears to be important to biological activity. By length is meant the distance between the A ring and the first charge bearing N of substituent Q. For example, when the A moiety is a 5 or 6-member ring [e.g. the benzene moiety of the benzodiazepinedione nucleus (carbon 7 of Formula II)], suitable lengths for bivalent radical L<sup>1</sup> range from about 3 to about 9 methylene equivalents. L<sup>1</sup> is therefore a bivalent radical
- 10        containing from 3 to 9 methylene groups connecting Q to the number 7 carbon of the benzodiazepinedione nucleus (or an equivalent position on other 5 or 6-member A rings) where any methylene group or groups may be replaced with one or more or a combination of; alkene, alkyne, aryl, heterocycle or a functional group or groups containing the heteroatoms N, O, and S, so long as the overall length is equivalent to from 3 to 9 methylene groups. These functional
- 15        groups include one or more of the following:

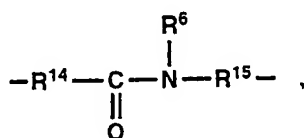


- and may be isolated within the linker (e.g. forming ethers, thioethers, ketones, sulfoxides and the like) or combined in any combination, provided only that the compounds so produced are stable in aqueous solution and do not exceed the above stated length requirements. For example, combining these functional groups produces esters, amides, ureidos, carbamates, carbonates, sulfonamides, sulfoxides, sulfones, and the like. Preferred lengths for L<sup>1</sup> are from 4 to 6 while most preferred lengths are about 5 methylene equivalents. In counting atoms
- 25        comprising L<sup>1</sup>, only those atoms sequentially linking Q with ring A are counted except when a homo- or heterocycle comprises L<sup>1</sup> in which case the fewest number of atoms separating these moieties are counted. Furthermore, given that a bivalent radical is of appropriate length, it is preferred that it be somewhat rigid, that is, contain one or more sp<sup>2</sup> or sp atoms.

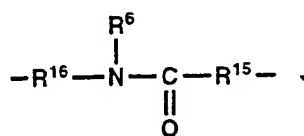
- In the description that follows, the free valence to the left of the page is bonded to Q, while the free valence to the right is bonded to carbon 7 of the compound of formula II (or an equivalent position on other 5 or 6-member A rings) Preferred L<sup>1</sup>'s are selected from
- 30

- substituted or unsubstituted; C3-C7-alkylene, C3-C7-cycloalkylene, C3-C7-alkenylene, C4-C7-cycloalkenylene, C5-C8-cycloalkadienylene, C3-C7-alkadienylene, C3-C7-alkynylene, C4-C7-alkenynylene, C6-C14-arylene, C6-C14-aryl-C2-C4-alkynylene, C1-C3-alkyl-C6-C14-aryl-C2-C4-alkynylene, C6-C14-aryl-C2-C4-alkenylene, C1-C3-alkyl-C6-C14-arylene, C1-C3-alkyl-C6-C14-aryl-C2-C4-alkenylene, C6-C14-aryl-C1-C3-alkylene, C6-C14-aryl-C1-C3-alkoxyene, C1-C3-alkyl-C6-C14-aryl-C1-C2-alkylene, C1-C3-alkoxy-C6-C14-arylene, C2-C8-alkoxyene, C1-C5-alkoxy-C1-C5-alkylene, C6-C10-aryloxyene, C6-C10-aryloxy-C1-C5-alkylene, C6-C10-arylthio-C1-C5-alkylene,

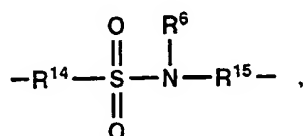
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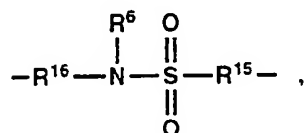
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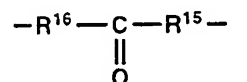
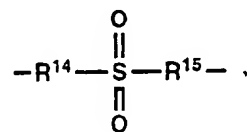
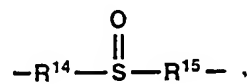
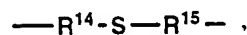
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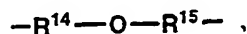
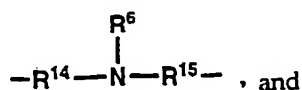


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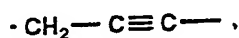




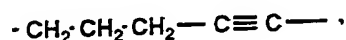
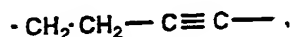
5 where  $\text{R}^{14}$  is selected from; a chemical bond, C<sub>1</sub>-C<sub>8</sub>-alkyl, C<sub>3</sub>-C<sub>7</sub>-cycloalkyl, C<sub>2</sub>-C<sub>5</sub>-alkenyl, C<sub>3</sub>-C<sub>5</sub>-alkynyl, C<sub>6</sub>-C<sub>10</sub>-aryl, C<sub>1</sub>-C<sub>3</sub>-alkyl-C<sub>6</sub>-C<sub>12</sub>-aryl, C<sub>1</sub>-C<sub>2</sub>-alkyl-C<sub>6</sub>-C<sub>10</sub>-aryl-C<sub>1</sub>-C<sub>2</sub>-alkyl, C<sub>6</sub>-C<sub>10</sub>-aryl-C<sub>1</sub>-C<sub>2</sub>-alkyl, and C<sub>6</sub>-C<sub>10</sub>-aryloxy-C<sub>1</sub>-C<sub>2</sub>-alkyl.  $\text{R}^{15}$  is selected from; a chemical bond, C<sub>1</sub>-C<sub>4</sub>-alkyl, C<sub>2</sub>-C<sub>4</sub>-alkenyl, C<sub>2</sub>-C<sub>4</sub>-alkynyl, C<sub>6</sub>-C<sub>10</sub>-aryl, and C<sub>1</sub>-C<sub>3</sub>-alkyl-C<sub>6</sub>-C<sub>12</sub>-aryl.  $\text{R}^{16}$  is  
10 selected from; a chemical bond, C<sub>1</sub>-C<sub>5</sub>-alkyl, C<sub>3</sub>-C<sub>7</sub>-cycloalkyl, C<sub>3</sub>-C<sub>5</sub>-alkenyl, C<sub>3</sub>-C<sub>5</sub>-alkynyl, C<sub>6</sub>-C<sub>10</sub>-aryl, C<sub>1</sub>-C<sub>3</sub>-alkyl-C<sub>6</sub>-C<sub>12</sub>-aryl, and C<sub>6</sub>-C<sub>10</sub>-aryl-C<sub>1</sub>-C<sub>2</sub>-alkyl. The substituents are preferably selected from one to three  $\text{R}^{12}$  defined above.

More preferred bivalent radicals  $\text{L}^1$  are selected from the following groups:

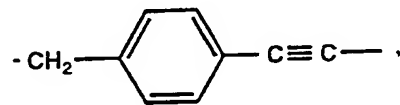
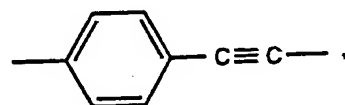
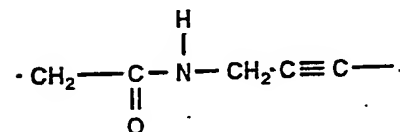
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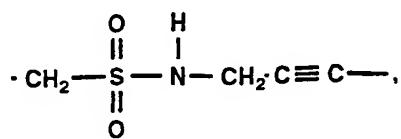
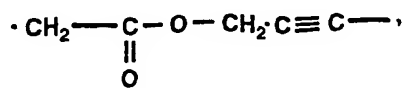
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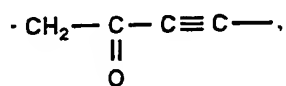
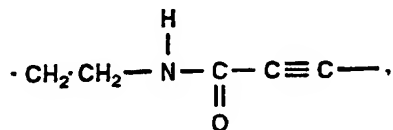


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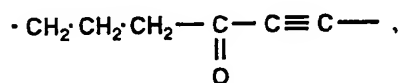
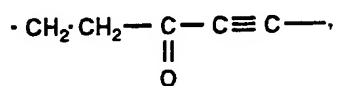
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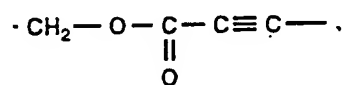
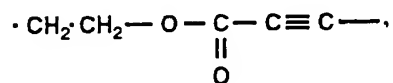
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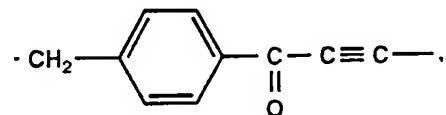
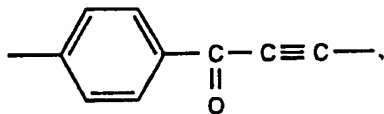
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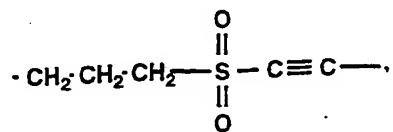
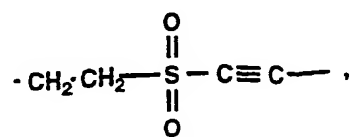
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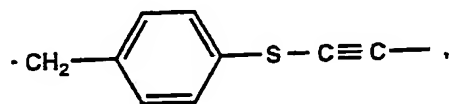
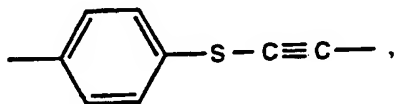
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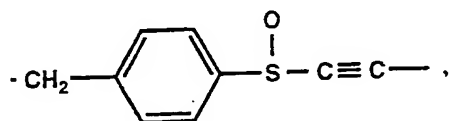
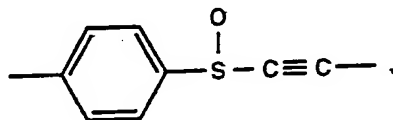
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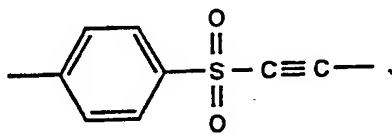
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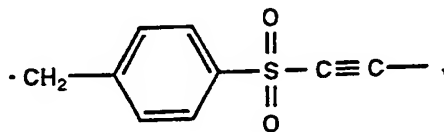
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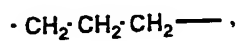
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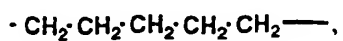
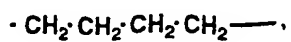


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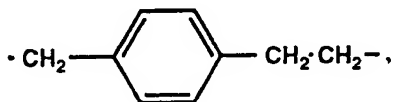
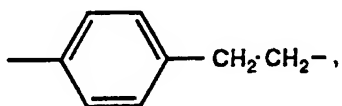
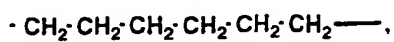
Group 3

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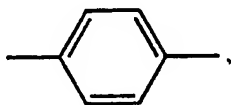


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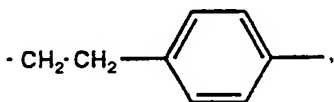
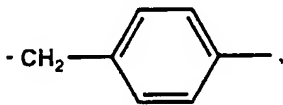


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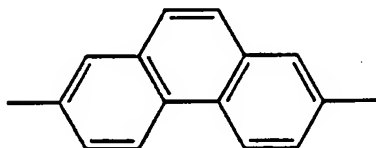
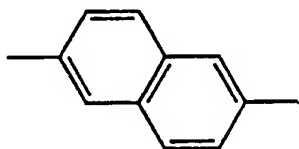
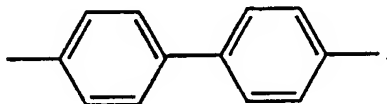
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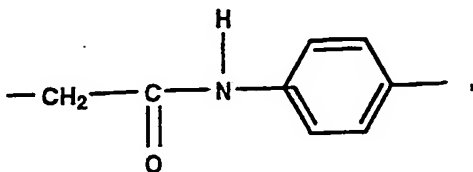
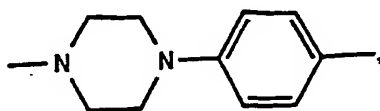
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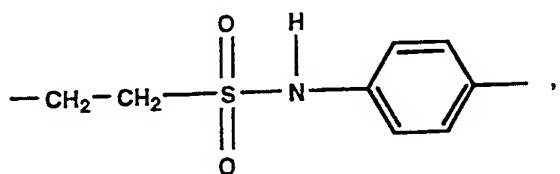
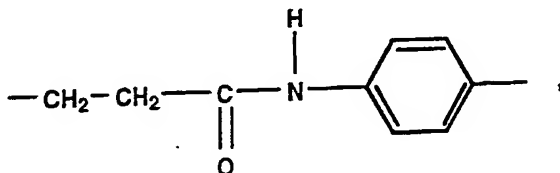
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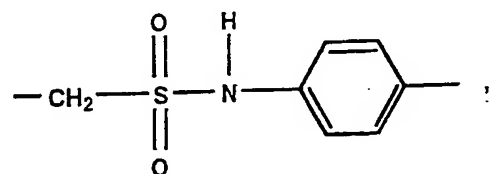
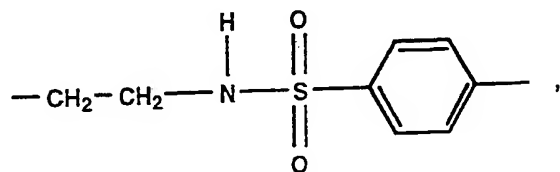
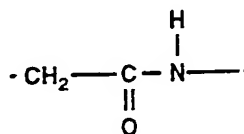
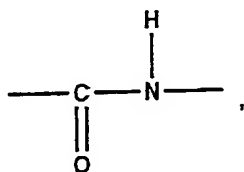
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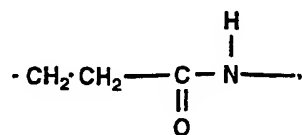
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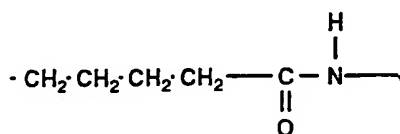
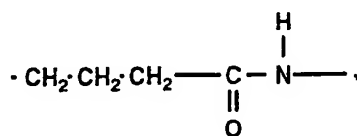
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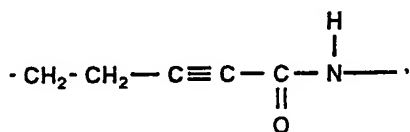
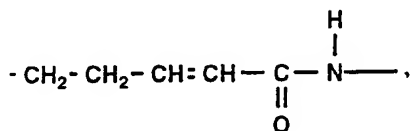




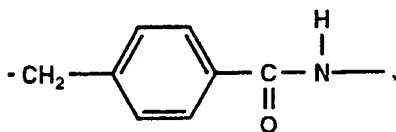
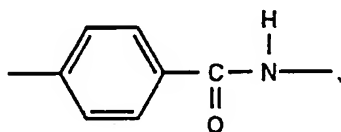
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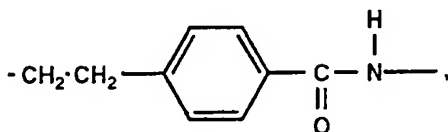
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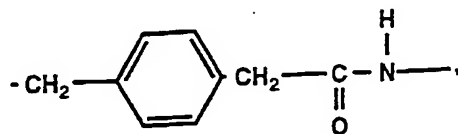
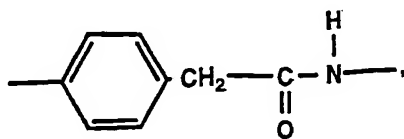


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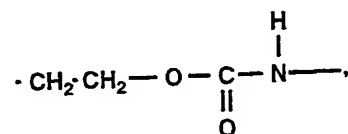


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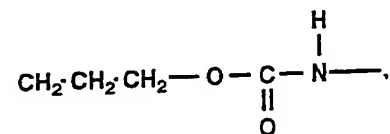




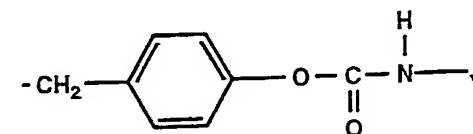
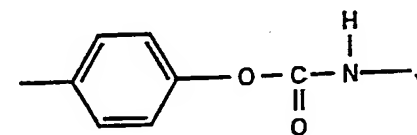
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Group 6

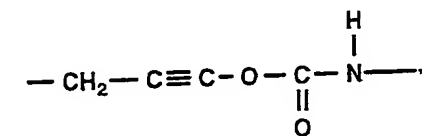
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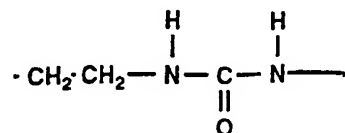


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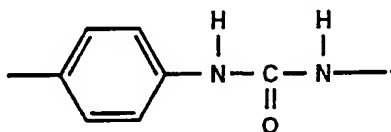
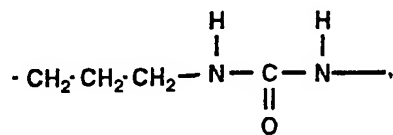


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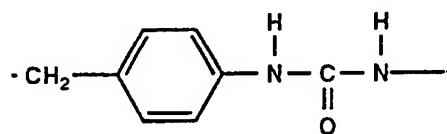


Group 7

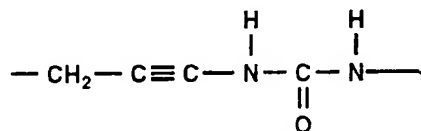
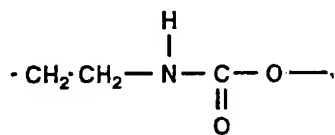
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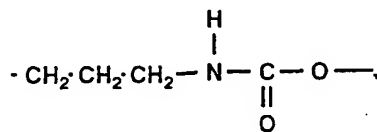
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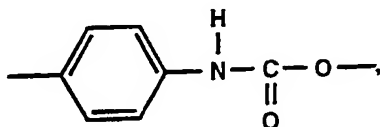
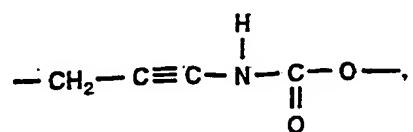


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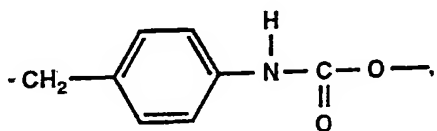
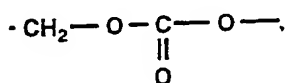
Group 8

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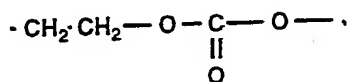




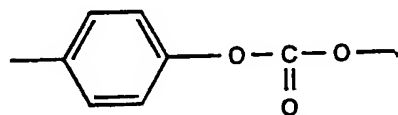
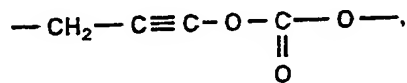
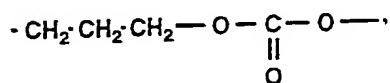
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Group 9

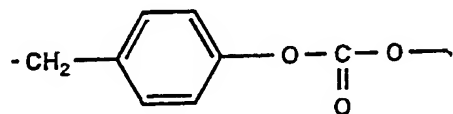
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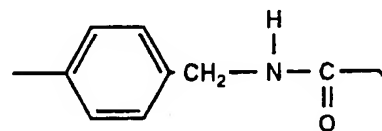
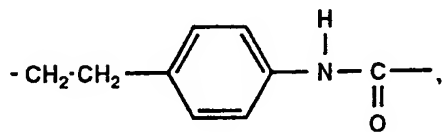
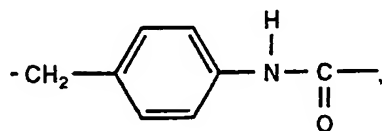
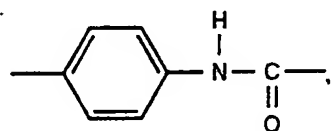
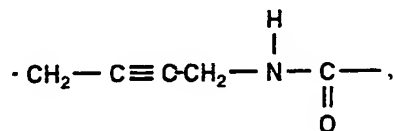
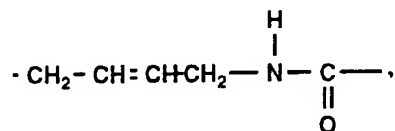
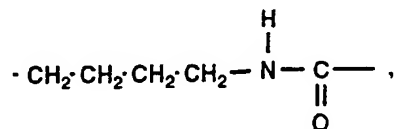
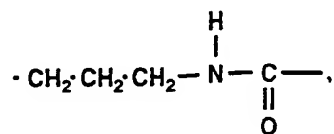
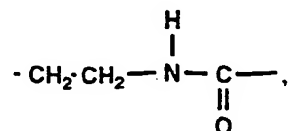


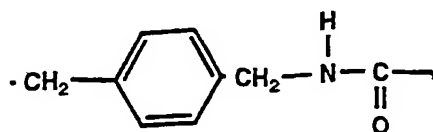
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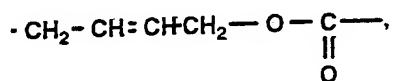
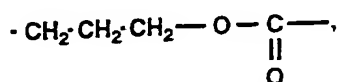
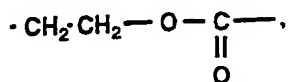
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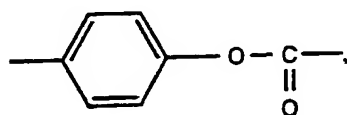
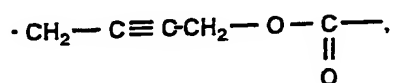
Group 10

Group 11

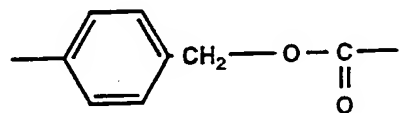
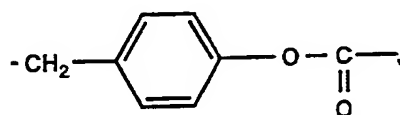
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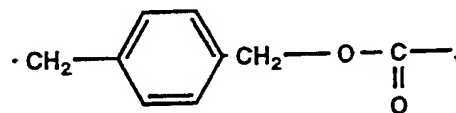
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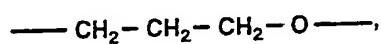
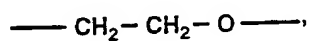
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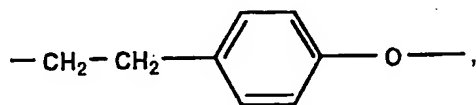
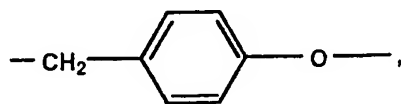
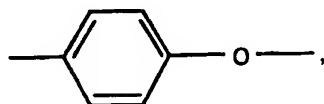
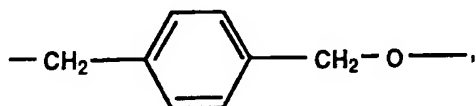
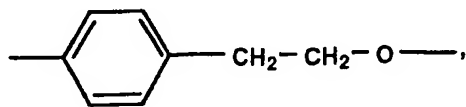
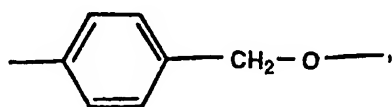


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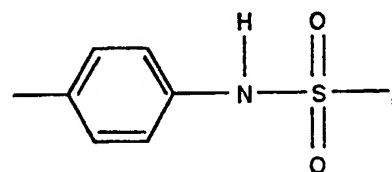
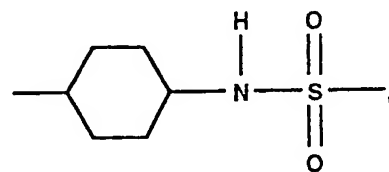
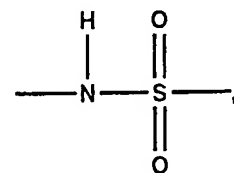
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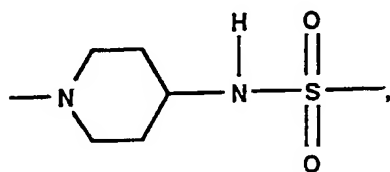
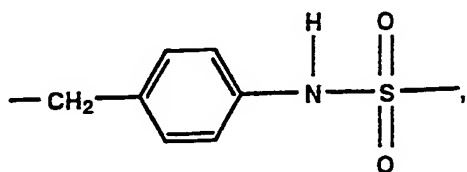
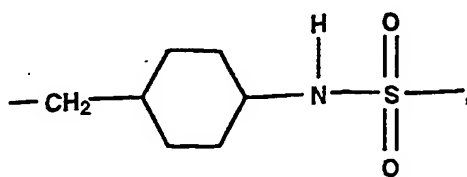
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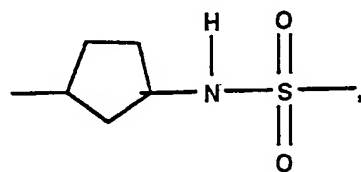
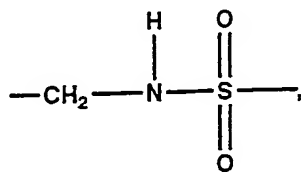


## Group 13

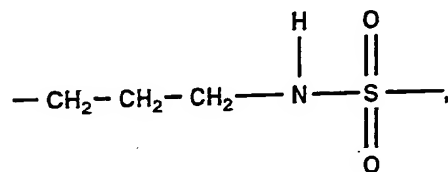
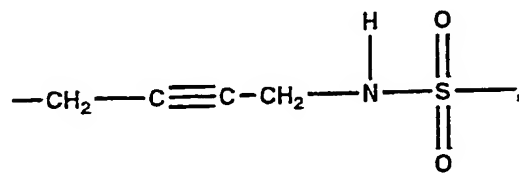
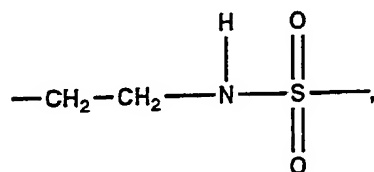




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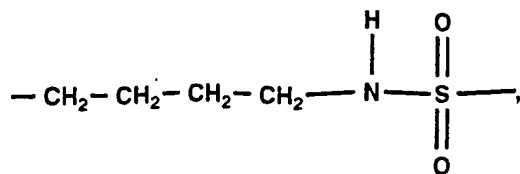


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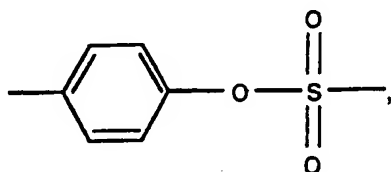


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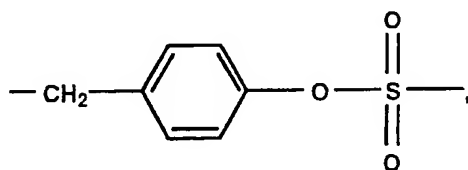
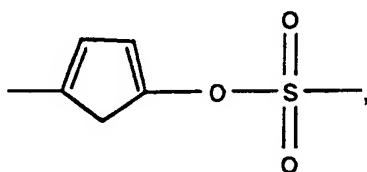




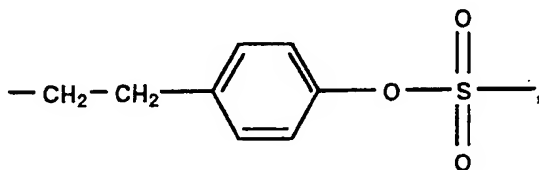
Group 14



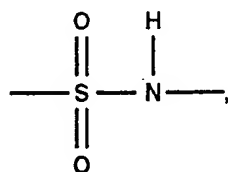
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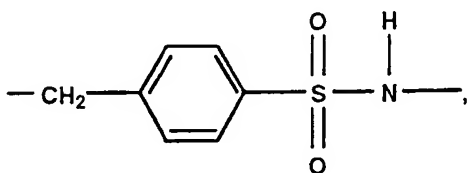
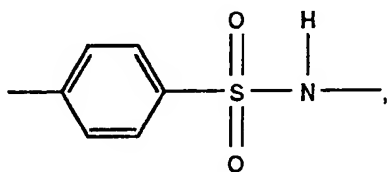
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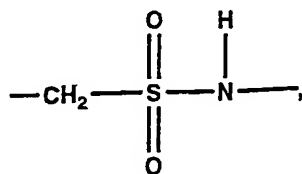
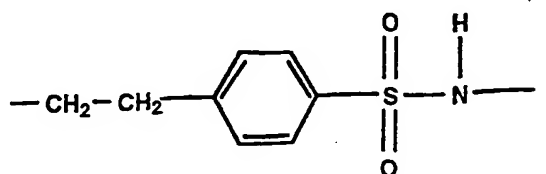


Group 15

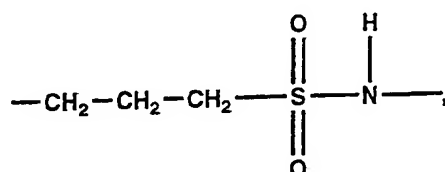
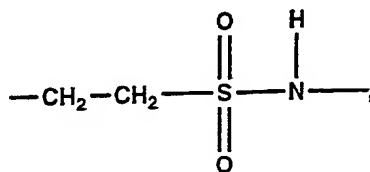


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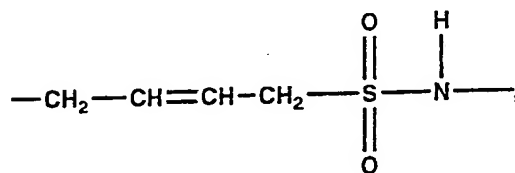
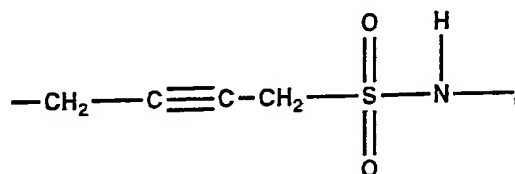
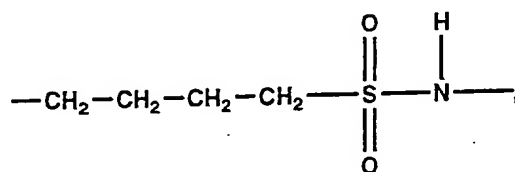




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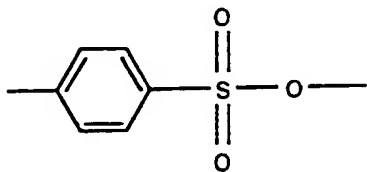


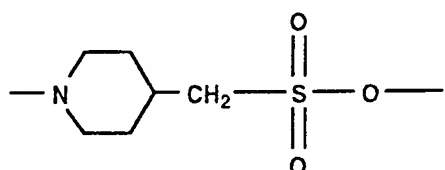
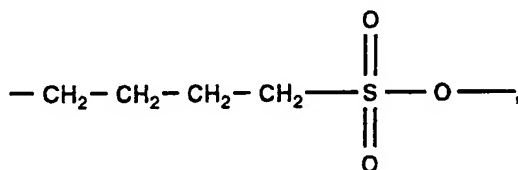
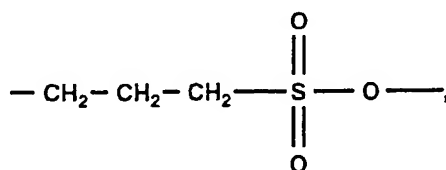
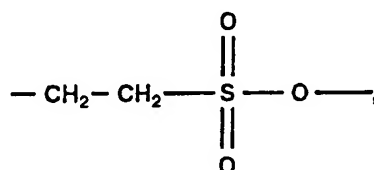
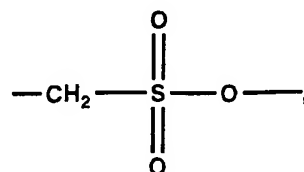
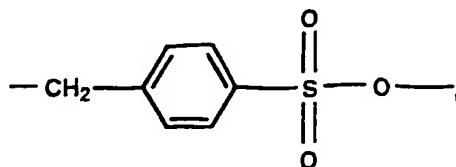
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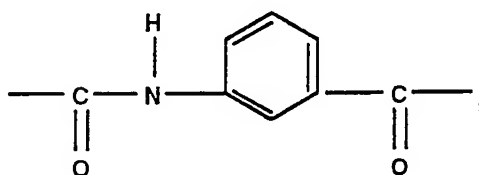
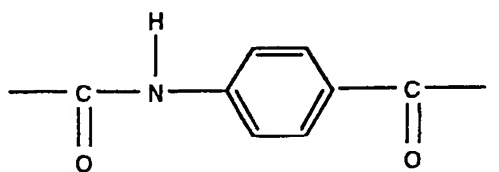
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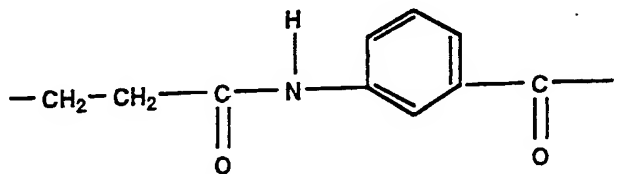
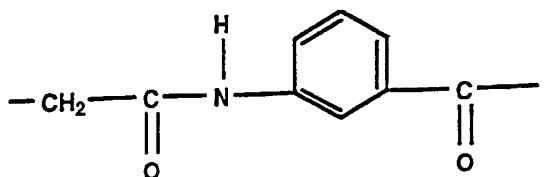
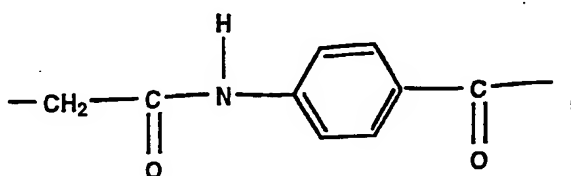
Group 16



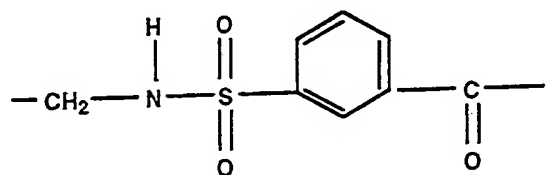
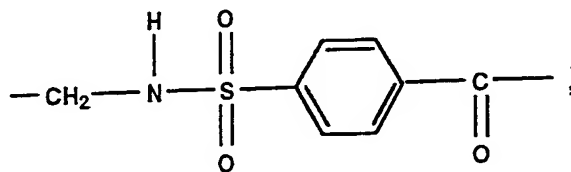


Group 17

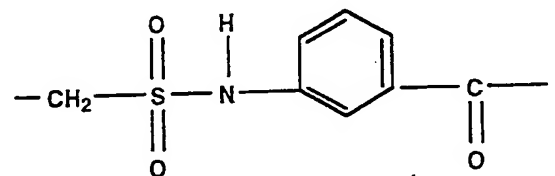
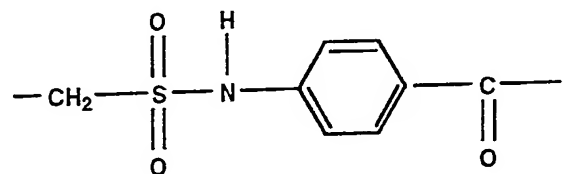




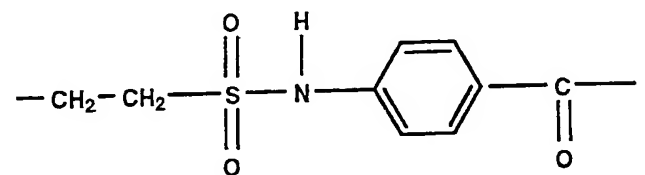
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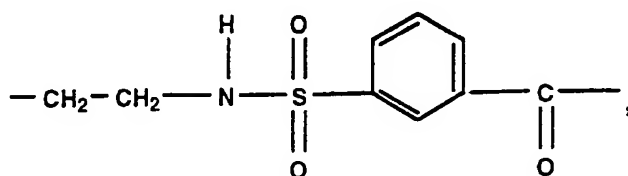
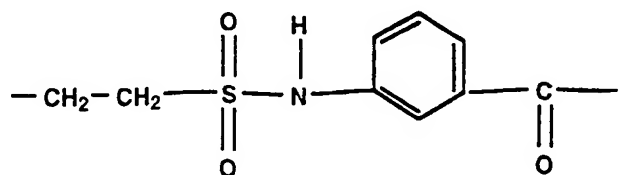


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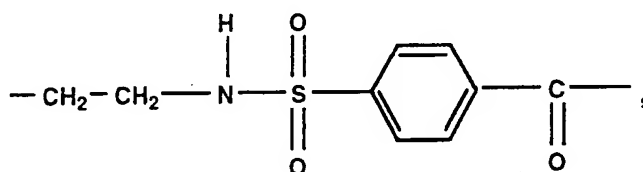


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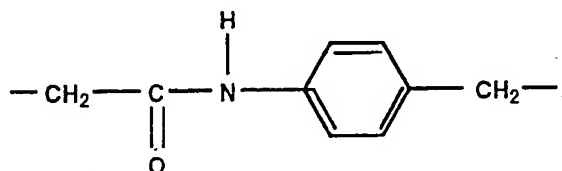
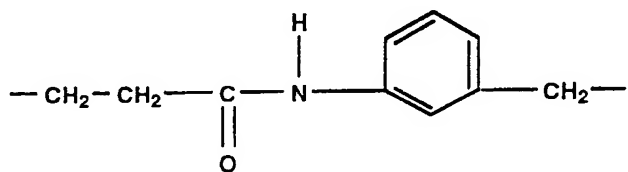




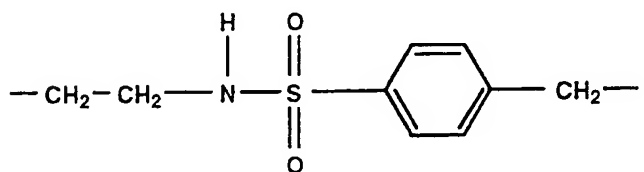
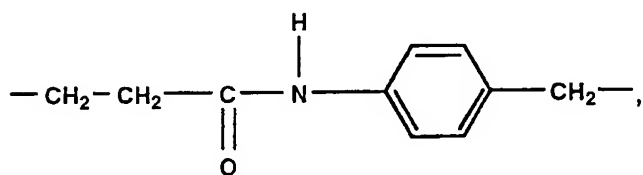
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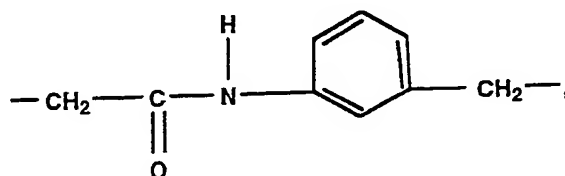
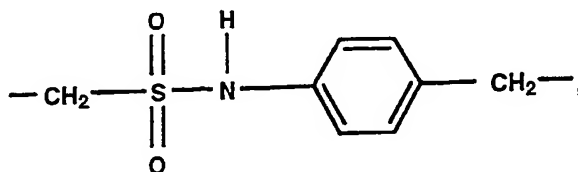
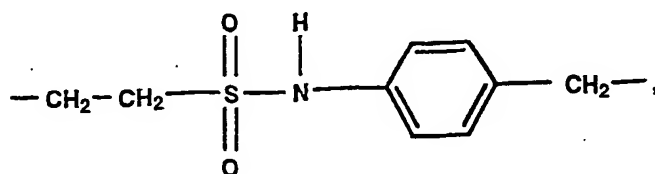
10 Group 18



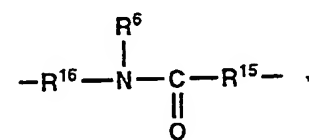
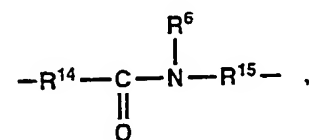
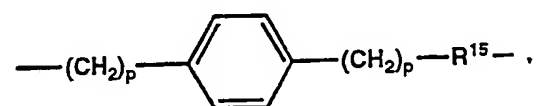
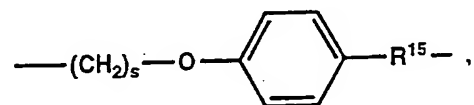
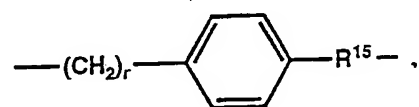
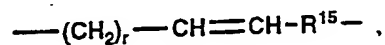
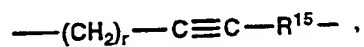
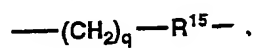
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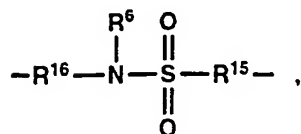
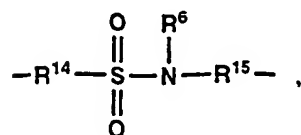


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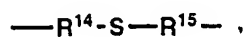


Most preferred L<sup>1</sup>'s are selected from;

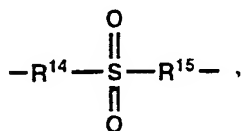
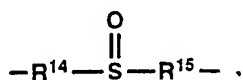




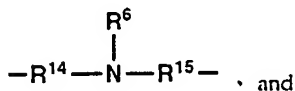
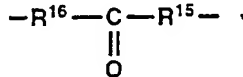
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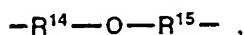
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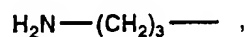


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where p is 1, 2, 3, or 4; q is 3, 4, 5, 6 or 7; r is 1, 2, 3, 4, or 5; s is 2 or 3. R<sup>14</sup> is selected from; a  
 25 chemical bond, C<sub>1</sub>-C<sub>5</sub>-alkyl, C<sub>3</sub>-C<sub>7</sub>-cycloalkyl, C<sub>2</sub>-C<sub>5</sub>-alkenyl, C<sub>3</sub>-C<sub>5</sub>-alkynyl, C<sub>6</sub>-C<sub>10</sub>-aryl, C<sub>1</sub>-  
 C<sub>2</sub>-alkyl-C<sub>6</sub>-C<sub>12</sub>-aryl, C<sub>1</sub>-C<sub>2</sub>-alkyl-C<sub>6</sub>-C<sub>10</sub>-aryl-C<sub>1</sub>-C<sub>2</sub>-alkyl, C<sub>6</sub>-C<sub>10</sub>-aryl-C<sub>1</sub>-C<sub>2</sub>-alkyl, C<sub>6</sub>-C<sub>10</sub>-  
 aryloxy-C<sub>1</sub>-C<sub>2</sub>-alkyl, and piperiziny. R<sup>15</sup> is a chemical bond connecting L<sup>1</sup> to position 7 of the  
 benzodiazepinedione. R<sup>16</sup> is selected from; C<sub>1</sub>-C<sub>5</sub>-alkyl, C<sub>3</sub>-C<sub>7</sub>-cycloalkyl, C<sub>3</sub>-C<sub>5</sub>-alkenyl, C<sub>3</sub>-  
 C<sub>5</sub>-alkynyl, C<sub>6</sub>-C<sub>10</sub>-aryl, C<sub>1</sub>-C<sub>3</sub>-alkyl-C<sub>6</sub>-C<sub>12</sub>-aryl, C<sub>6</sub>-C<sub>10</sub>-aryl-C<sub>1</sub>-C<sub>2</sub>-alkyl, and piperiziny.  
 30 For the foregoing most preferred L<sup>1</sup>'s R<sup>14</sup> and R<sup>16</sup> bond L<sup>1</sup> to Q<sup>1</sup>.

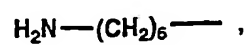
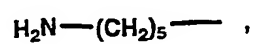
Most preferred Q<sup>1</sup>-L<sup>1</sup>- combinations are selected from;



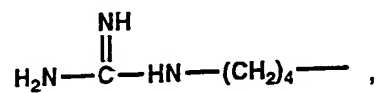
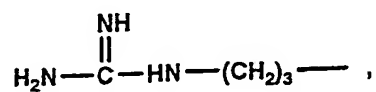
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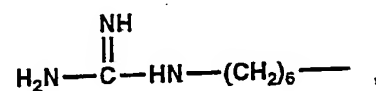
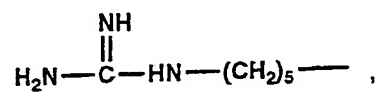
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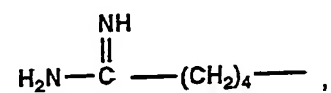
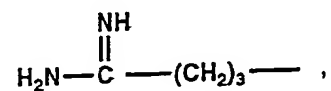
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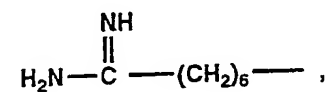
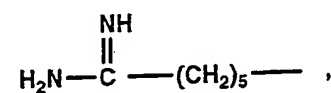
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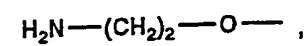
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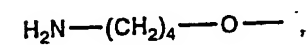
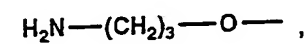
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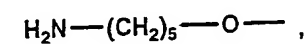
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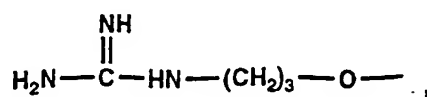
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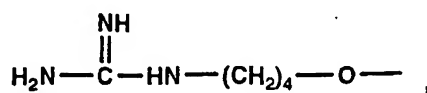
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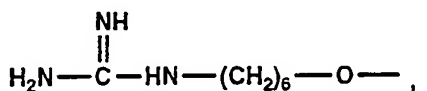
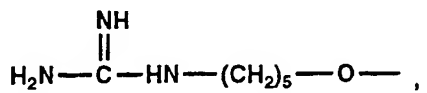




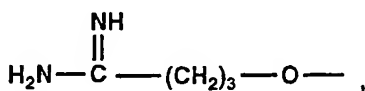
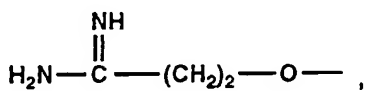
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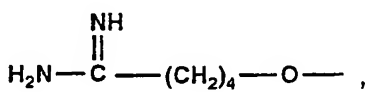
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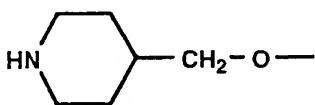
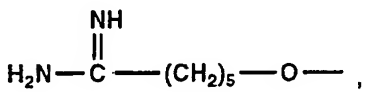
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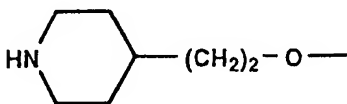
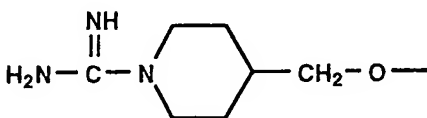
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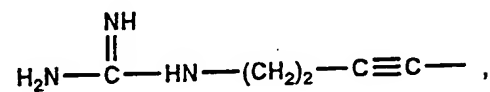
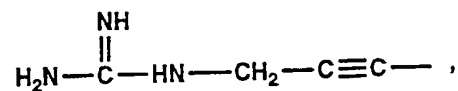
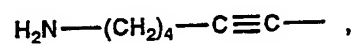
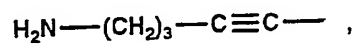
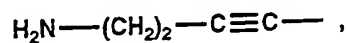
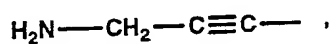
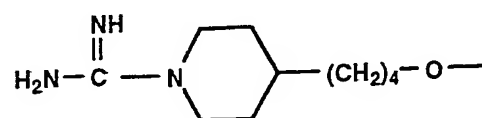
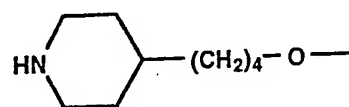
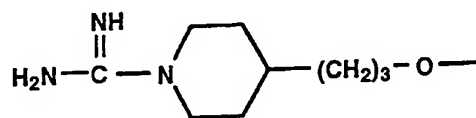
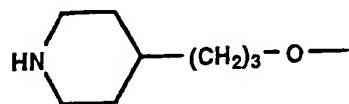
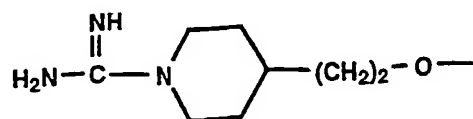


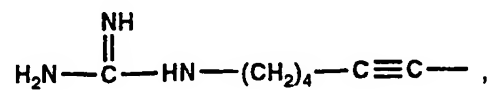
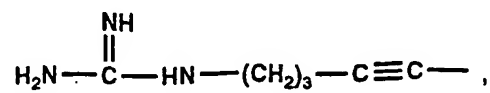
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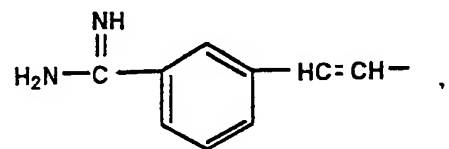
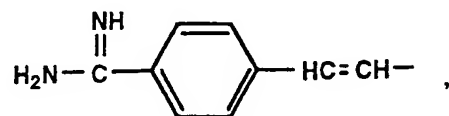
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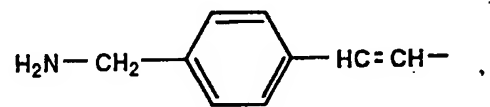




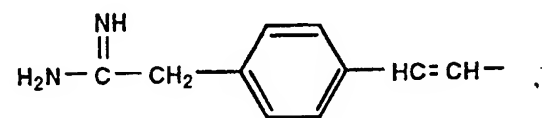
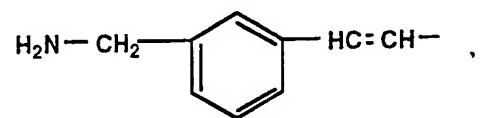
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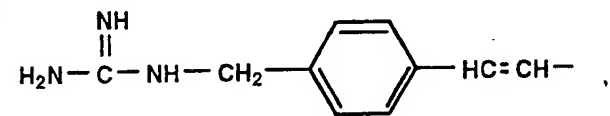
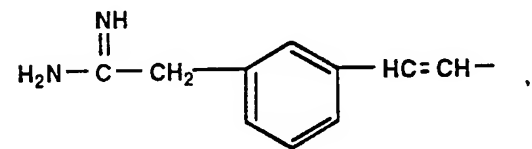
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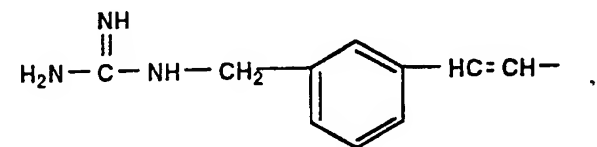
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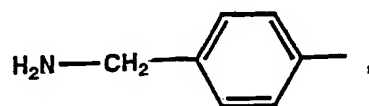


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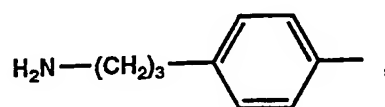
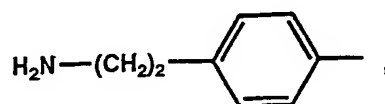


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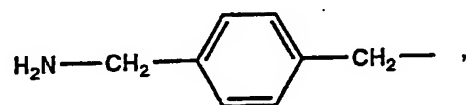
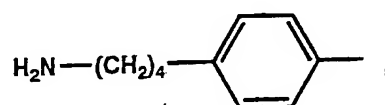




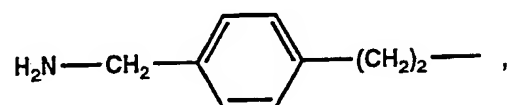
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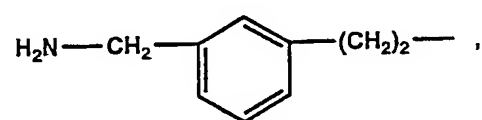
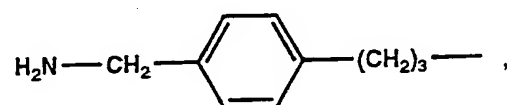
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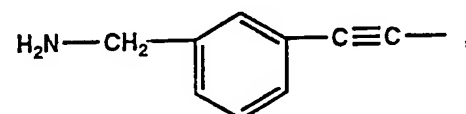
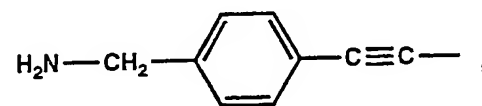
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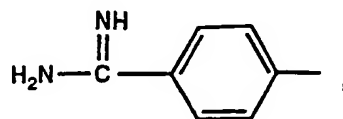
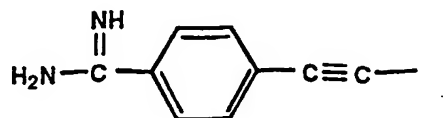
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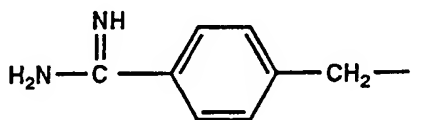
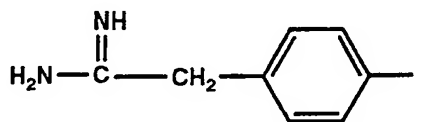
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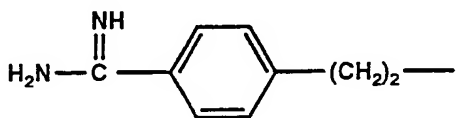
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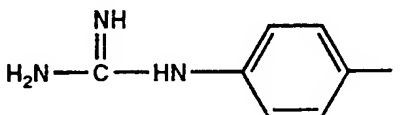
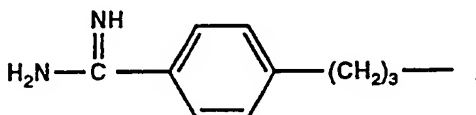
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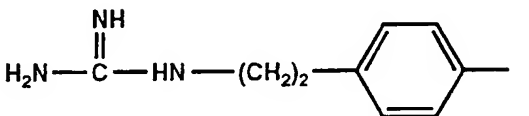
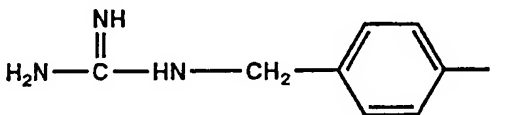
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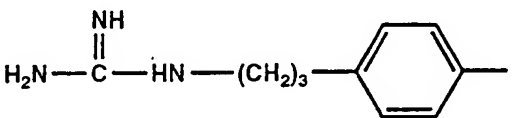
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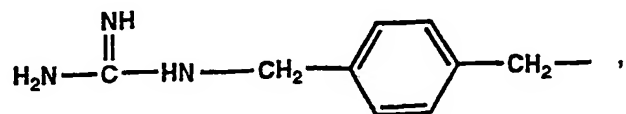
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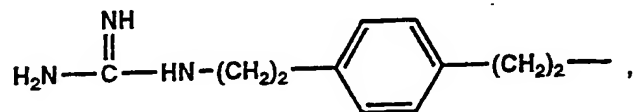
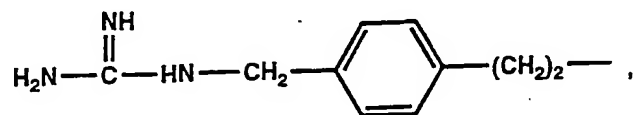
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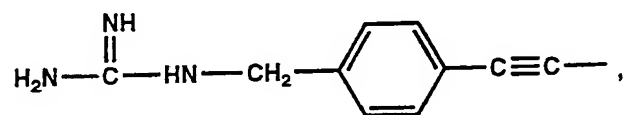
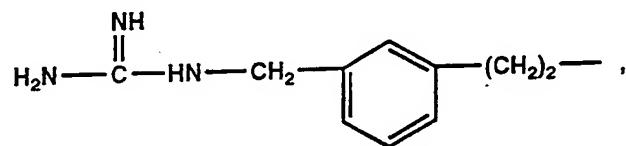
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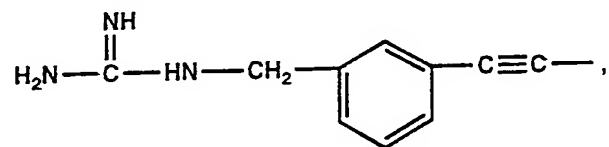
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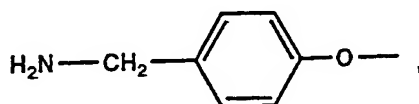
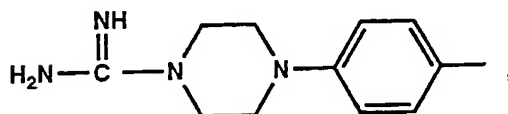
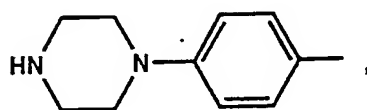
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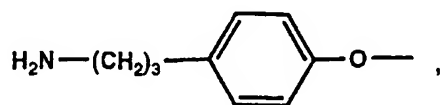
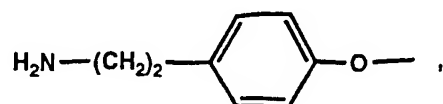


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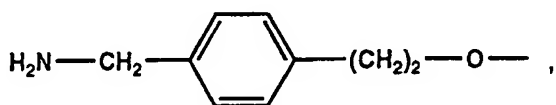
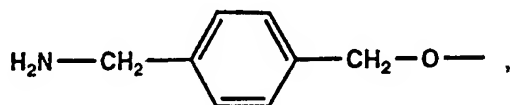


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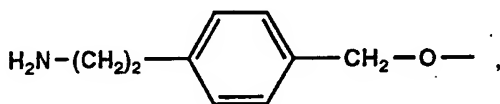




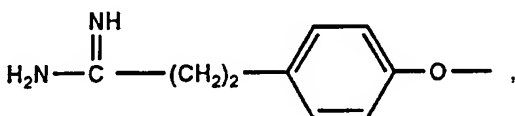
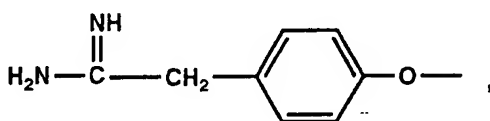
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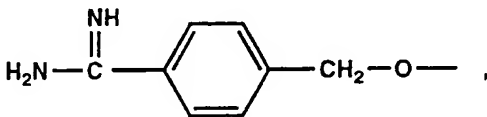
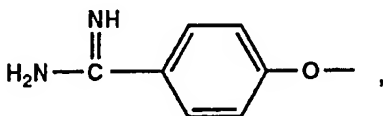
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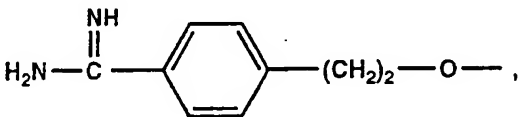
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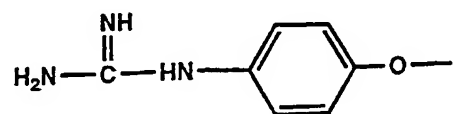
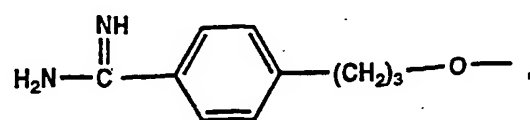
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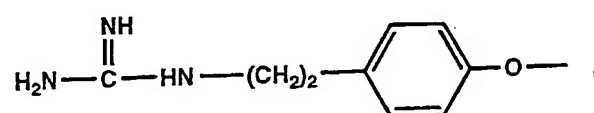
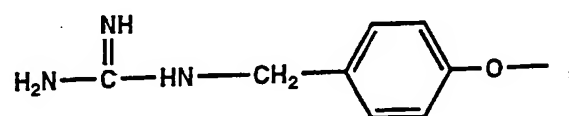
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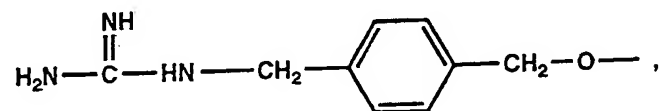
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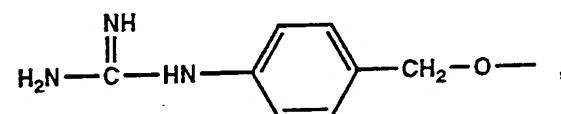
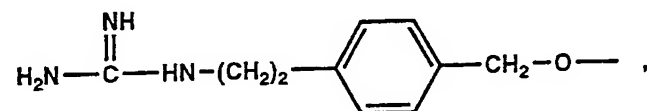
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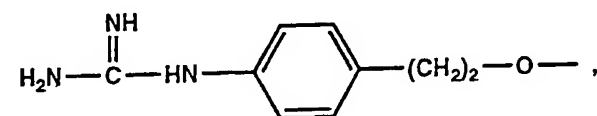
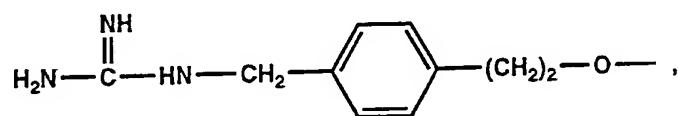
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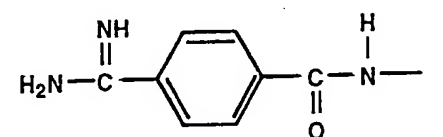
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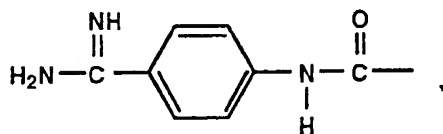
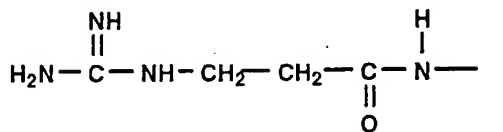
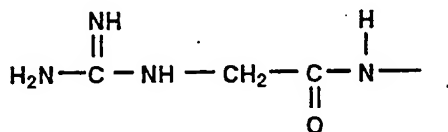
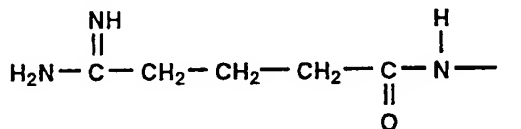
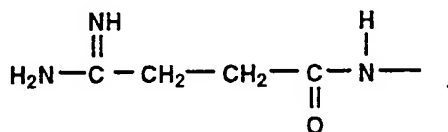
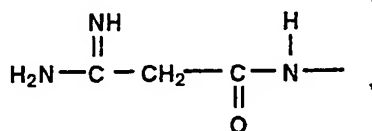
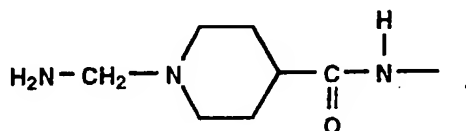
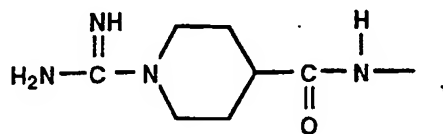
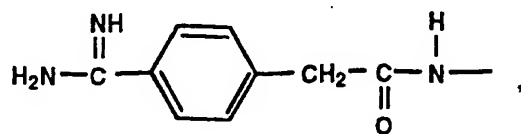
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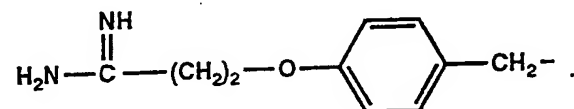
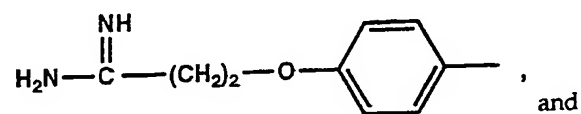
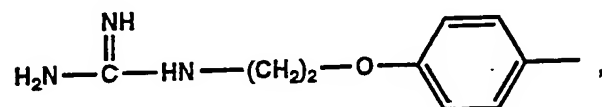
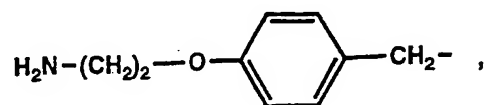
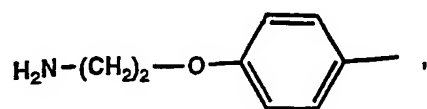


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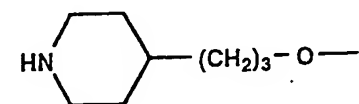
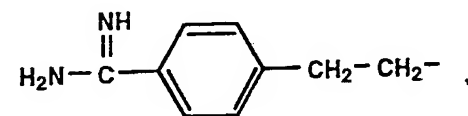
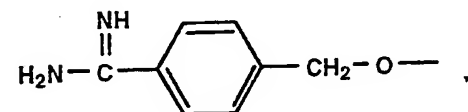
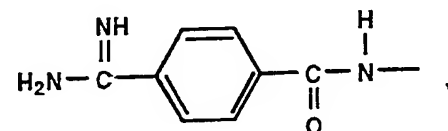


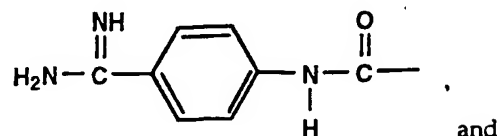
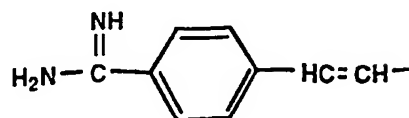




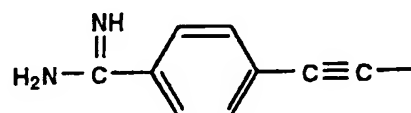


Very most preferred Q<sup>1</sup>-L<sup>1</sup> combinations are selected from;





and

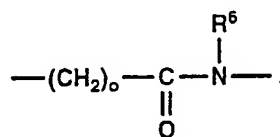


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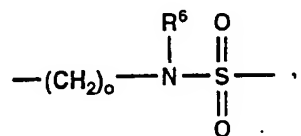
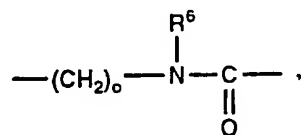
The length of the bivalent radical  $\text{L}^2$  is also important to biological activity. Here, when the A moiety is a 10-member fused ring [e.g. the naphthalene moiety of the naphthadiazepinedione nucleus (carbon 8 of Formula III)], suitable lengths for bivalent radical  $\text{L}^2$  range from about 1 to about 7 methylene equivalents.  $\text{L}^2$  is therefore a bivalent radical containing from 1 to 7 methylene groups connecting  $\text{Q}^1$  to the number 8 carbon of the naphthadiazepinedione nucleus (or an equivalent position on other 10-member A rings) where any methylene group or groups may be replaced with one or more alkene, alkyne, aryl, heterocycle or functional groups containing the heteroatoms N, O, and S, so long as the overall length is equivalent to 1 to 7 methylene groups. Here, the functional groups are the same as described above for  $\text{L}^1$  and may be combined in any combination provided again that the compounds produced are stable in aqueous solution and do not exceed the stated length requirements for  $\text{L}^2$ . Preferred lengths for  $\text{L}^2$  are from 2 to 4 while most preferred lengths are about 3 methylene equivalents.

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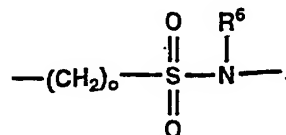
Preferred  $\text{L}^2$ 's are optionally substituted bivalent radicals selected from;



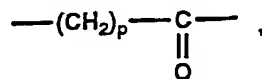
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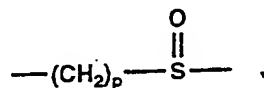
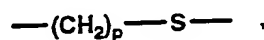
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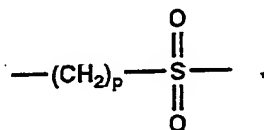
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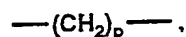
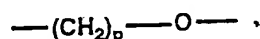
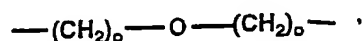
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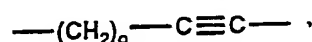
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where o is 0, 1 or 2, p is 1, 2, 3 or 4,  $R^6$  is defined above, and where the substituents are selected from one to three  $R^{12}$  groups defined above.

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#### c. Ring A

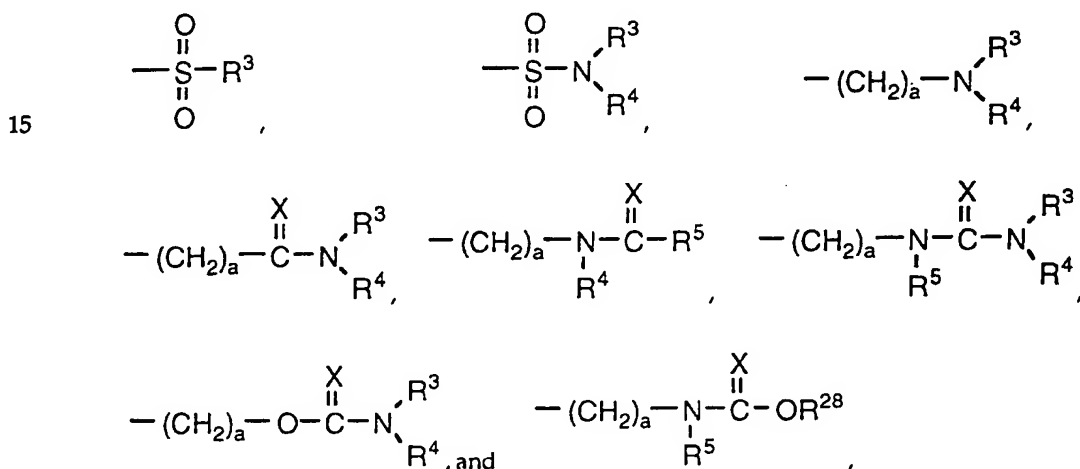
Preferred A rings include; benzene, naphthalene, tetrahydronaphthalene, cyclohexane, thiophene, pyridine, pyrazine, pyrimidine, quinoline, isoquinoline, quinoxaline, quinazoline, pteridine, and naphthyridine. The most preferred A ring is benzene.

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Preferred substituents of rings A are  $R^1$  and  $R^2$ , where each  $R^1$  and  $R^2$ , is independently selected from hydrogen, halogen(F, Cl, Br, I), cyano, nitro, carboxyl, mercaptocarbonyl, mercapto, phthalimido and a substituent having from 1 to 12 carbon atoms selected from: alkyl, hydroxyalkyl, formyl, haloalkyl, cycloalkyl, cycloalkylalkyl, cycloalkylhaloalkyl, cycloalkylcarbonyl, alkoxy, aralkyl, aralkylhaloalkyl, aryl, aroyl, aryloxy, aryloxyalkyl, aralkoxy, alkoxyalkyl, alkylcarbonyl, alkylcarbonylalkyl, alkoxycarbonyl,

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alkenyl, cycloalkenyl, alkylthio, carboxylalkyl, alkylcarbonyloxy, alkylcarbonyloxyalkyl, alkoxyalkyl, aralkoxyalkyl, aralkylcarbonyloxyalkyl, mercaptothiocarbonyl, mercaptoalkyl, alkoxyalkyl, alkylthio, cycloalkylthio, cycloalkylalkylthio, alkylthiocarbonyl, alkylcarbonylthio, alkylthiocarbonyloxy, alkylthiocarbonylthio, 5 alkylthiothiocarbonyl, alkylthiothiocarbonylthio, arylthio, arylthiocarbonyl, arylcarbonylthio, arylthiocarbonyloxy, arylthiocarbonylthio, arylthiothiocarbonyl, arylthiothiocarbonylthio, aralkylthio, aralkylthiocarbonyl, aralkylcarbonylthio, aralkylthiocarbonyloxy, aralkylthiocarbonylthio, alkylthiocarbonyl, aralkylthiocarbonylthio, alkylsulfanyl, alkylsulfonyl, aralkylsulfanyl, aralkylsulfonyl, arylsulfanyl, arylsulfonyl, pthalimidoalkyl, heteroaryl, 10 heteroarylalkyl, cycloheteroalkyl, cycloheteroalkylalkyl, and cycloheteroalkylcarbonylalkyl where each heteroaryl- and cyclohetero-containing groups has one or more ring hetero atoms selected from oxygen, sulfur and nitrogen. Optionally, each  $R^1$  and  $R^2$  also may be independently selected from sulfonyl, sulfonylamido, amino and amido radicals of the formula



where X is oxygen or sulfur, each "a" is a number independently selected from zero to six, inclusive, and where each of  $R^3$  through  $R^5$  is independently selected from the groups described above. Optionally,  $R^3$  and  $R^4$  taken together,  $R^4$  and  $R^5$  taken together and  $R^3$  and  $R^5$  taken together may each form a heterocyclic having from five to seven ring members including the hetero atom of the sulfonyl, amino or amido radical and which heterocyclic may 25 further contain one or more hetero atoms as ring members selected from oxygen, nitrogen, and sulfur and which heterocyclic group may be saturated or partially saturated.  $R^{28}$  is selected from the groups defined below.

30 d. The 7-member "lactam" ring containing TUG;

Preferred seven member rings are those of formula I where T bears the substituent  $R^{22}$ . Thus T is preferably  $\text{sp}^3$  carbon or nitrogen. Also preferably G is a methylene group. The most preferred "lactam ring" is a diazepinedione.

## e. Substituents of T-U-G

R<sup>19</sup> and R<sup>20</sup> are preferably hydrogen or halogen. R<sup>18</sup> and R<sup>21</sup> are preferably hydrogen or halogen or taken together form oxo. R<sup>22</sup> is preferably selected from; hydrogen, optionally substituted C<sub>1</sub>-C<sub>6</sub>-alkyl, where the substituents are selected from amino, hydroxy, halo (F, Cl, Br, I), carboxy, and C<sub>1</sub>-C<sub>4</sub>-alkoxycarbonyl, optionally substituted C<sub>6</sub>-C<sub>10</sub>-aryl, and optionally substituted C<sub>6</sub>-C<sub>12</sub>-aryl-C<sub>1</sub>-C<sub>4</sub>-alkyl, where the substituents on any aryl group are selected from amino, nitro, halo (F, Cl, Br, I), halo(F, Cl, Br, I)-C<sub>1</sub>-C<sub>6</sub>-alkyl, C<sub>1</sub>-C<sub>6</sub>-alkoxy, C<sub>6</sub>-C<sub>10</sub>-aryloxy, amino-C<sub>1</sub>-C<sub>6</sub>-acylamino, amino-C<sub>1</sub>-C<sub>6</sub>-acyl, and guanidino C<sub>6</sub>-C<sub>10</sub>-arylamino.

## f. The amino acid linking moiety D

D is preferably hydrogen, phenyl, or lower alkyl. Optionally, D is -(C=O)-Xaa, where Xaa is one to three natural, unnatural, or modified  $\alpha$ -amino acid residues, preferably one or two naturally occurring amino acid residues, and most preferably a naturally occurring hydrophobic amino acid residue. Optionally, when Xaa is two amino acid residues, the terminal residue is Arg or Lys amide. Common naturally occurring  $\alpha$ -amino acids are described by the standard three letter amino acid code when referring to amino acids or residues. When the three-letter code begins with a lower-case letter, it is understood the amino acid is the unnatural or D-isomeric form. Standard abbreviations are listed in The Merck Index, 10th Edition, pp Misc-2 - Misc-3. Modified or unusual  $\alpha$ -amino acids such as norleucine (Nle) and ornithine (Orn) are designated as described in U.S. Patent and Trademark Office Official Gazette 1114TMOG, May 15, 1990.

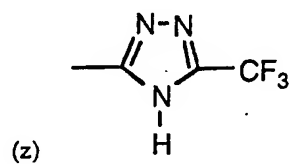
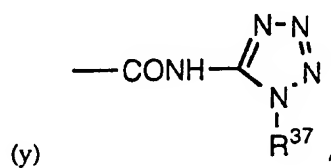
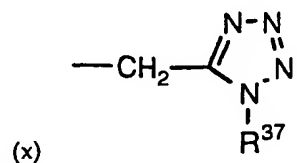
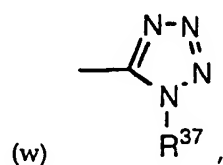
## g. The negatively charged acidic W moiety

W is -R<sup>27</sup>-w.

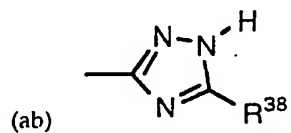
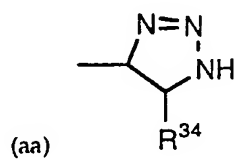
R<sup>27</sup> is preferably selected from (a) a covalent bond, (b) substituted or unsubstituted methylene, and (c) substituted or unsubstituted ethylene, where the substituents are independently selected from one or more of the groups (i) nitro, (ii) halo(F, Cl, Br, I), (iii) C<sub>1</sub>-C<sub>6</sub>alkyl, (iv) halo(F, Cl, Br, I)-C<sub>1</sub>-C<sub>6</sub>alkyl, and (v) substituted or unsubstituted phenyl where the substituents are selected from (1) C<sub>1</sub>-C<sub>6</sub>alkyl, (2) C<sub>1</sub>-C<sub>6</sub>alkoxy, (3) halo(F, Cl, Br, I), and (4) CF<sub>3</sub>.

"w" is selected from (a) -COR<sup>28</sup>, (b) -SO<sub>3</sub>R<sup>31</sup>, (c) -NHSO<sub>2</sub>R<sup>32</sup>, (d) -PO(OR<sup>31</sup>)<sub>2</sub>, (e) -SO<sub>2</sub>NHR<sup>32</sup>, (f) -CONHOR<sup>31</sup>, (g) -C(OH)R<sup>33</sup>PO(OR<sup>33</sup>)<sub>2</sub>, (h) -CN, (i) -SO<sub>2</sub>NH-heterocycle where the heterocycle is a 5- or 6-member aromatic ring containing 1 to 3 heteroatoms selected from O, N, and S and where the heterocycle is unsubstituted or substituted with one or two substituents selected from the group (i) -OH, (ii) -SH, (iii) -(C<sub>1</sub>-C<sub>4</sub>alkyl), (iv) -(C<sub>1</sub>-C<sub>4</sub>alkoxyl), (v) CF<sub>3</sub>, (vi) halo(F, Cl, Br, I), (vii) NO<sub>2</sub>, (viii) -COOH, (ix) -COO-(C<sub>1</sub>-C<sub>4</sub>alkyl), (x) -NH<sub>2</sub>, (xi) -NH(C<sub>1</sub>-C<sub>4</sub>alkyl), or (xii) -N(C<sub>1</sub>-C<sub>4</sub>alkyl)<sub>2</sub>, (j) -CH<sub>2</sub>SO<sub>2</sub>NH-heterocycle, (k) -SO<sub>2</sub>NHCOR<sup>33</sup>, (l) -CH<sub>2</sub>SO<sub>2</sub>NHCOR<sup>32</sup>, (m) -CONHSO<sub>2</sub>R<sup>33</sup>, (n) -CH<sub>2</sub>CONHSO<sub>2</sub>R<sup>33</sup>, (o) -NHCONHSO<sub>2</sub>R<sup>33</sup>, (p) -NHSO<sub>2</sub>NHCOR<sup>33</sup>, (q) -CONHNHSO<sub>2</sub>CF<sub>3</sub>, (r) CON(OH)R<sup>31</sup>, (s) -CONHCOCF<sub>3</sub>, (t) -CONHSO<sub>2</sub>R<sup>28</sup>, (u) -CONHSO<sub>2</sub>R<sup>29</sup>, (v) -CONHSO<sub>2</sub>R<sup>30</sup>,

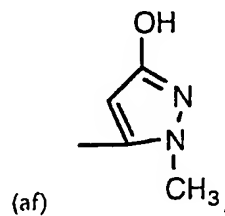
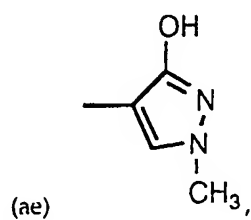
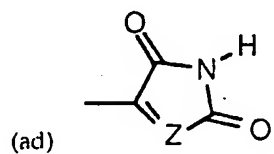
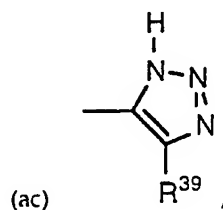
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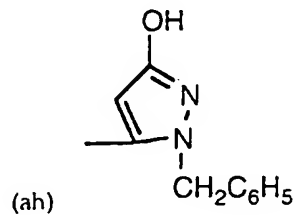
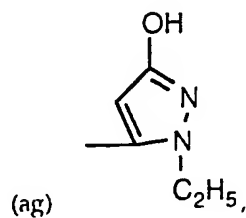
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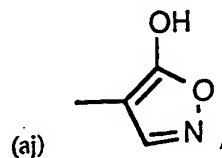
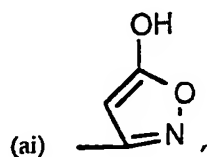


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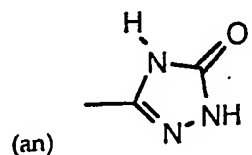
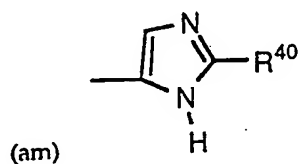
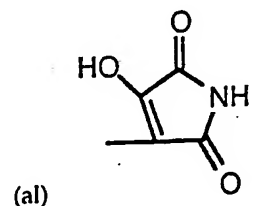
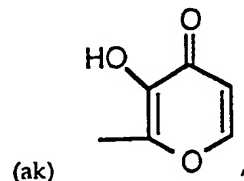


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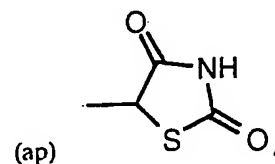
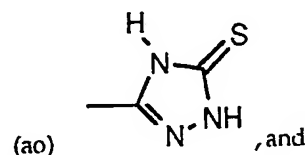




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15  $R^{28}$  is selected from the group consisting of (a) hydroxy, (b) C<sub>1</sub>-C<sub>8</sub>-alkoxy, (c) C<sub>3</sub>-C<sub>12</sub>-alkenoxy, (d) C<sub>6</sub>-C<sub>12</sub>-aryloxy, (e) C<sub>1</sub>-C<sub>6</sub>-alkyl-C<sub>6</sub>-C<sub>12</sub>-aryloxy, (f) di-C<sub>1</sub>-C<sub>8</sub>-alkylamino-C<sub>1</sub>-C<sub>8</sub>-alkoxy, (g) acylamino-C<sub>1</sub>-C<sub>8</sub>-alkoxy selected from the group (i) acetylaminoethoxy, (ii) nicotinoylaminoethoxy, and (iii) succinamidoethoxy, (h) C<sub>1</sub>-C<sub>8</sub>-alkoxyloxy-C<sub>1</sub>-C<sub>8</sub>-alkoxy, (i) C<sub>6</sub>-C<sub>12</sub>-aryl-C<sub>1</sub>-C<sub>8</sub>-alkoxy where the aryl group is unsubstituted or substituted with one to three  
20 of the groups (i) nitro, (ii) halo (F, Cl, Br, I), (iii) C<sub>1</sub>-C<sub>4</sub>-alkoxy, and (iv) amino, (j) hydroxy-C<sub>2</sub>-C<sub>8</sub>-alkoxy, (k) dihydroxy-C<sub>3</sub>-C<sub>8</sub>-alkoxy, and (l)  $NR^{29}R^{30}$ .

$R^{29}$  and  $R^{30}$  are independently selected from the group (a) hydrogen, (b) C<sub>1</sub>-C<sub>8</sub>-alkyl, (c) C<sub>3</sub>-C<sub>8</sub>-alkenyl, (d) C<sub>6</sub>-C<sub>12</sub>-aryl where the aryl group is unsubstituted or substituted with one to three of the groups (i) nitro, (ii) halo (F, Cl, Br, I), (iii) C<sub>1</sub>-C<sub>4</sub>-alkoxy, and (iv) amino, and  
25 (e) C<sub>6</sub>-C<sub>12</sub>-aryl-C<sub>1</sub>-C<sub>8</sub>-alkyl where the aryl group is unsubstituted or substituted with one to three of the groups (i) nitro, (ii) halo (F, Cl, Br, I), and (iii) C<sub>1</sub>-C<sub>4</sub>-alkoxy.

$R^{31}$  is selected from the group consisting of (a) H, (b) C<sub>1</sub>-C<sub>6</sub> alkyl, (c) halo(F, Cl, Br, I)-C<sub>1</sub>-C<sub>6</sub> alkyl, (d) phenyl, (e) benzyl, and (f) -CH<sub>2</sub>-O-COCH<sub>3</sub>.



$R^{32}$  is selected from the group consisting of (a) H, (b) benzyl, and (c)  $-\text{CH}(\text{R}^{35})-\text{O}-\text{C}(\text{O})\text{R}^{35}$ .

$R^{33}$  is selected from the group consisting of (a) aryl, (b) heterocycle, (c)  $(\text{C}_3-\text{C}_7)$ -cycloalkyl, (d)  $(\text{C}_1-\text{C}_4)$ -alkyl, unsubstituted or substituted with a substituent selected from the group consisting of (i) aryl, (ii) heterocycle, (iii)  $-\text{OH}$ , (iv)  $-\text{SH}$ , (v)  $(\text{C}_1-\text{C}_4)$ -alkyl, (vi)  $(\text{C}_1-\text{C}_4)$ -alkoxy, (vii)  $(\text{C}_1-\text{C}_4)$ -alkylthio, (viii)  $-\text{CF}_3$ , (ix) halo (F, Cl, Br, I), (x)  $-\text{NO}_2$ , (xi)  $-\text{CO}_2\text{H}$ , (xii)  $\text{CO}_2$ - $(\text{C}_1-\text{C}_4)$ -alkyl, (xiii)  $-\text{NH}_2$ , (xiv)  $-\text{N}[(\text{C}_1-\text{C}_4)\text{-alkyl}]_2$ , (xv)  $-\text{NH}[(\text{C}_1-\text{C}_4)\text{-alkyl}]$ , (xvi)  $-\text{PO}_3\text{H}$ , or (xvii)  $\text{PO}(\text{OH})(\text{C}_1-\text{C}_4)\text{-alkoxy}$ , or (e)  $(\text{C}_1-\text{C}_4)$ -perfluoroalkyl.

$R^{34}$  is selected from the group consisting of (a)  $-\text{CN}$ , (b)  $-\text{NO}_2$ , (c)  $-\text{COOR}^{31}$ , (d)  $\text{C}_1-\text{C}_6$ -perfluoroalkyl, and (e)  $\text{CF}_3$ .

$R^{35}$  is independently selected from the group consisting of (a) H, (b)  $(\text{C}_1-\text{C}_6)$ -alkyl,  $(\text{C}_2-\text{C}_6)$ -alkenyl,  $(\text{C}_2-\text{C}_6)$ -alkynyl or  $(\text{C}_3-\text{C}_8)$ -cycloalkyl, each of which is unsubstituted or substituted with: (i) OH, (ii)  $(\text{C}_1-\text{C}_4)$ -alkoxy, (iii)  $\text{CO}_2\text{R}^{33}$ , (iv)  $\text{OCOR}^{33}$ , (v)  $\text{CONHR}^{33}$ , (vi)  $\text{CON}(\text{R}^{33})_2$ , (vii)  $\text{N}(\text{R}^{33})\text{C}(\text{O})\text{R}^{33}$ , (viii)  $\text{NH}_2$ , (ix)  $(\text{C}_1-\text{C}_4)$ -alkylamino, (x)  $\text{di}[(\text{C}_1-\text{C}_4)\text{-alkyl}]\text{amino}$ , (xi) aryl, (xii) heteroaryl, (c)  $-\text{C}(\text{O})\text{-aryl}$ , (d)  $-\text{NO}_2$ , (e) halo (Cl, Br, I, F), (f)  $-\text{OH}$ , (g)  $-\text{OR}^{36}$ , (h)  $(\text{C}_1-\text{C}_4)$ -perfluoroalkyl, (i)  $-\text{SH}$ , (j)  $-\text{S}(\text{O})_{1-2}(\text{C}_1-\text{C}_4)\text{-alkyl}$ , (k)  $\text{CO}_2\text{R}^{33}$ , (l)  $-\text{SO}_3\text{H}$ , (m)  $-\text{NR}^{33}\text{R}^{36}$ , (n)  $-\text{NR}^{33}\text{C}(\text{O})\text{R}^{36}$ , (o)  $-\text{NR}^{33}\text{COOR}^{32}$ , (p)  $-\text{SO}_2\text{NHR}^{32}$ , (q)  $-\text{SO}_2\text{NR}^{33}\text{R}^{33}$ , (r)  $-\text{NHSO}_2\text{R}^{32}$ , (s)  $-\text{C}(\text{O})\text{NHSO}_2\text{R}^{32}$ , (t) aryl, (u) heteroaryl, (v) morpholin-4-yl, (w)  $\text{CONH}_2$ , or (y) 1H-tetrazol-5-yl.

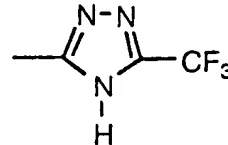
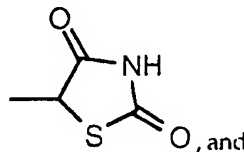
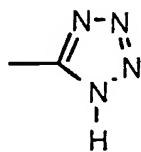
$R^{36}$  is selected from the group (a) H; or (b)  $(\text{C}_1-\text{C}_4)$ -alkyl unsubstituted or substituted with (i)  $\text{NH}_2$ , (ii)  $\text{NH}[(\text{C}_1-\text{C}_4)\text{-alkyl}]$ , (iii)  $\text{N}[(\text{C}_1-\text{C}_4)\text{-alkyl}]_2$ , (iv)  $\text{CO}_2\text{H}$ , (v)  $\text{CO}_2(\text{C}_1-\text{C}_4)\text{-alkyl}$ , (vi) OH, (vii)  $\text{SO}_3\text{H}$ , or (viii)  $\text{SO}_2\text{NH}_2$ .

$R^{37}$  is selected from the group consisting of (a) H, (b)  $(\text{C}_1-\text{C}_6)$ -alkyl, (c)  $(\text{C}_2-\text{C}_6)$ -alkenyl, (d)  $(\text{C}_1-\text{C}_6)$ -alkoxyalkyl, (e)  $-\text{CH}_2-\text{O}-\text{COCH}_3$ , or (f)  $-\text{CH}_2$ -phenyl, where the phenyl is unsubstituted or substituted with a substituent selected from  $-\text{NO}_2$ ,  $-\text{NH}_2$ ,  $-\text{OH}$ , or  $-\text{OCH}_3$ .

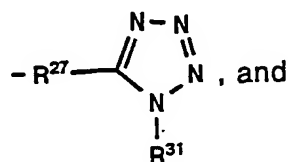
$R^{38}$ ,  $R^{39}$ , and  $R^{40}$  are each independently selected from H, Cl, CN,  $\text{NO}_2$ ,  $\text{CF}_3$ ,  $\text{C}_2\text{F}_5$ ,  $\text{C}_3\text{F}_7$ ,  $\text{CHF}_2$ ,  $\text{CH}_2\text{F}$ ,  $\text{CO}_2\text{CH}_3$ ,  $\text{CO}_2\text{C}_2\text{H}_5$ ,  $\text{SO}_2\text{CH}_3$ ,  $\text{SO}_2\text{CF}_3$  and  $\text{SO}_2\text{C}_6\text{F}_5$ ; wherein Z is selected from O, S,  $\text{NR}^{41}$  and  $\text{CH}_2$ .

$R^{41}$  is selected from hydrogen,  $\text{CH}_3$ , and  $\text{CH}_2\text{C}_6\text{H}_5$  or a tautomer or pharmaceutically acceptable salt thereof.

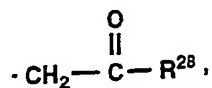
Preferred "w"'s are selected from:  $-\text{COR}^{28}$ ,  $-\text{NHSO}_2\text{CF}_3$ .



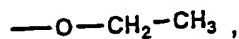
Most preferred W's are either a tetrazole or a carboxylate. Preferably these groups are represented by;



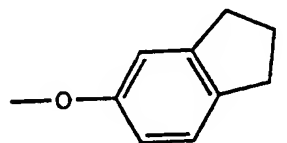
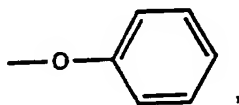
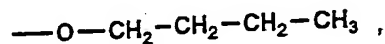
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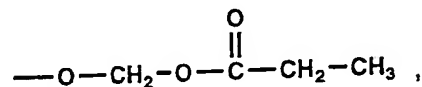
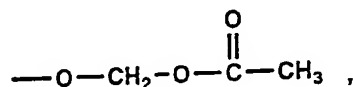
where  $R^{27}$  is methylene, and  $R^{28}$  and  $R^{31}$  are hydroxy and hydrogen respectively. Optionally,  $R^{28}$  and  $R^{31}$  are substituents that transform compounds represented by structural formulae I-  
 10 VI into prodrugs. When W contains a carboxylate, for example, preferred prodrug forms include simple esters,  $\alpha$ -acyloxyalkyl esters, and amides. Preferred esters include compounds where  $R^{28}$  is selected from;



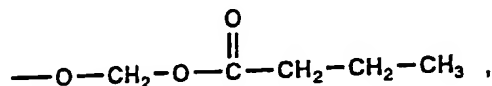
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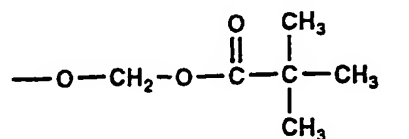
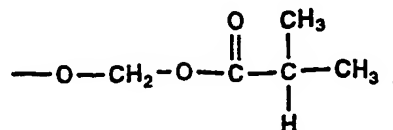


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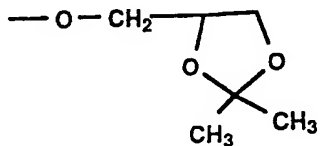


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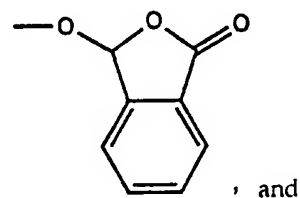
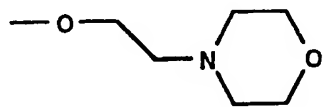




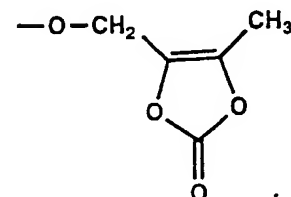
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The most preferred compounds of this invention are represented by structural formula

II and are selected from the following:

- 1-(4-chlorophenyl)methyl-4-(2-carboxyethyl)-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-  
 20 benzodiazapine-2,5-dione.  
 1-(diphenyl)methyl-4-(2-carboxyethyl)-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-  
 benzodiazapine-2,5-dione.

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- 1-(*m*-trifluoromethyl)phenyl-4-(2-carboxyethyl)-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-methyl-4-(2-carboxyethyl)-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 5 1-(1-methyl)ethyl-4-(2-carboxyethyl)-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(2-methyl)propyl-4-(2-carboxyethyl)-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(4-chlorophenyl)methyl-4-(2-carboxyethyl)-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 10 1-(diphenyl)methyl-4-(2-carboxyethyl)-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(*m*-trifluoromethyl)phenyl-4-(2-carboxyethyl)-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 15 1-methyl-4-(2-carboxyethyl)-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(1-methyl)ethyl-4-(2-carboxyethyl)-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(2-methyl)propyl-4-(2-carboxyethyl)-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 20 1-(diphenyl)methyl-4-(2-carboxyethyl)-7-(6-aminoethyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(diphenyl)methyl-4-(2-carboxyethyl)-7-(6-aminoethyl-1-ynyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 25 1-(diphenyl)methyl-4-(2-carboxyethyl)-7-(5-aminopentoxyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(diphenyl)methyl-4-(2-carboxyethyl)-7-(5-guanidinopentyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(diphenyl)methyl-4-(2-carboxyethyl)-7-(4-guanidinobutoxy)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 30 1-(diphenyl)methyl-4-(2-carboxyethyl)-7-(5-aminopent-1-ynyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(diphenyl)methyl-4-(2-carboxyethyl)-7-(5-guanidinopent-1-ynyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 35 1-(diphenyl)methyl-4-(2-carboxyethyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(diphenyl)methyl-4-(2-carboxyethyl)-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.

- 1-(diphenyl)methyl-4-(2-carboxyethyl)-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(diphenyl)methyl-4-(2-carboxyethyl)-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 5 1-(4-chlorophenyl)methyl-4-(2-carboxyethyl)-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(4-chlorophenyl)methyl-4-(2-carboxyethyl)-7-(5-aminopentoxyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(4-chlorophenyl)methyl-4-(2-carboxyethyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 10 1-(4-chlorophenyl)methyl-4-(2-carboxyethyl)-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(4-chlorophenyl)methyl-4-(2-carboxyethyl)-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 15 1-(4-chlorophenyl)methyl-4-(2-carboxyethyl)-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-methyl-4-(2-carboxyethyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-methyl-4-(2-carboxyethyl)-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 20 1-methyl-4-(2-carboxyethyl)-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-methyl-4-(2-carboxyethyl)-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 25 1-(1-methyl)ethyl-4-(2-carboxyethyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(1-methyl)ethyl-4-(2-carboxyethyl)-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(1-methyl)ethyl-4-(2-carboxyethyl)-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 30 1-(2-methyl)ethyl-4-(2-carboxyethyl)-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(2-methyl)propyl-4-(2-carboxyethyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 35 1-(2-methyl)propyl-4-(2-carboxyethyl)-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(2-methyl)propyl-4-(2-carboxyethyl)-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.

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- 1-(2-methyl)propyl-4-(2-carboxyethyl)-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(2,4-difluorophenyl)methyl-4-(2-carboxyethyl)-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 5 1-(2,4-difluorophenyl)methyl-4-(2-carboxyethyl)-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(2,4-difluorophenyl)methyl-4-(2-carboxyethyl)-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(2,4-difluorophenyl)methyl-4-(2-carboxyethyl)-7-(5-aminopentoxyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 10 1-(2,4-difluorophenyl)methyl-4-(2-carboxyethyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(2,4-difluorophenyl)methyl-4-(2-carboxyethyl)-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 15 1-(2,4-difluorophenyl)methyl-4-(2-carboxyethyl)-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(2,4-difluorophenyl)methyl-4-(2-carboxyethyl)-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(4-chlorophenyl)methyl-4-(2-carboxyethyl)-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 20 1-(diphenyl)methyl-4-(2-carboxyethyl)-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(*m*-trifluoromethyl)phenyl-4-(2-carboxyethyl)-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 25 1-methyl-4-(2-carboxyethyl)-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(1-methyl)ethyl-4-(2-carboxyethyl)-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(2-methyl)propyl-4-(2-carboxyethyl)-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 30 1-(2,4-difluorophenyl)methyl-4-(2-carboxyethyl)-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(4-chlorophenyl)methyl-4-(3-butanoyl)-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 35 1-(diphenyl)methyl-4-(3-butanoyl)-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(*m*-trifluoromethyl)phenyl-4-(3-butanoyl)-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.

- 1-methyl-4-(3-butanoyl)-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(1-methyl)ethyl-4-(3-butanoyl)-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 5 1-(2-methyl)propyl-4-(3-butanoyl)-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(4-chlorophenyl)methyl-4-(3-butanoyl)-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(diphenyl)methyl-4-(3-butanoyl)-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 10 1-(*m*-trifluoromethyl)phenyl-4-(3-butanoyl)-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-methyl-4-(3-butanoyl)-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 15 1-(1-methyl)ethyl-4-(3-butanoyl)-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(2-methyl)propyl-4-(3-butanoyl)-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(diphenyl)methyl-4-(3-butanoyl)-7-(6-aminoethyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 20 1-(diphenyl)methyl-4-(3-butanoyl)-7-(6-aminohex-1-ynyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(diphenyl)methyl-4-(3-butanoyl)-7-(5-aminopentoxyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 25 1-(diphenyl)methyl-4-(3-butanoyl)-7-(5-guanidinopentyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(diphenyl)methyl-4-(3-butanoyl)-7-(4-guanidinobutoxy)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(diphenyl)methyl-4-(3-butanoyl)-7-(5-aminopent-1-ynyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 30 1-(diphenyl)methyl-4-(3-butanoyl)-7-(5-guanidinopent-1-ynyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(diphenyl)methyl-4-(3-butanoyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 35 1-(diphenyl)methyl-4-(3-butanoyl)-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(diphenyl)methyl-4-(3-butanoyl)-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.

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- 1-(diphenyl)methyl-4-(3-butanoyl)-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(4-chlorophenyl)methyl-4-(3-butanoyl)-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 5 1-(4-chlorophenyl)methyl-4-(3-butanoyl)-7-(5-aminopentoxyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(4-chlorophenyl)methyl-4-(3-butanoyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(4-chlorophenyl)methyl-4-(3-butanoyl)-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 10 1-(4-chlorophenyl)methyl-4-(3-butanoyl)-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(4-chlorophenyl)methyl-4-(3-butanoyl)-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 15 1-methyl-4-(3-butanoyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-methyl-4-(3-butanoyl)-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-methyl-4-(3-butanoyl)-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 20 1-methyl-4-(3-butanoyl)-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(1-methyl)ethyl-4-(3-butanoyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(1-methyl)ethyl-4-(3-butanoyl)-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 25 1-(1-methyl)ethyl-4-(3-butanoyl)-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(2-methyl)ethyl-4-(3-butanoyl)-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 30 1-(2-methyl)propyl-4-(3-butanoyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(2-methyl)propyl-4-(3-butanoyl)-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(2-methyl)propyl-4-(3-butanoyl)-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 35 1-(2-methyl)propyl-4-(3-butanoyl)-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(2,4-difluorophenyl)methyl-4-(3-butanoyl)-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.



- 1-(2,4-difluorophenyl)methyl-4-(3-butanoyl)-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(2,4-difluorophenyl)methyl-4-(3-butanoyl)-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 5 1-(2,4-difluorophenyl)methyl-4-(3-butanoyl)-7-(5-aminopentoxyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(2,4-difluorophenyl)methyl-4-(3-butanoyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(2,4-difluorophenyl)methyl-4-(3-butanoyl)-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 10 1-(2,4-difluorophenyl)methyl-4-(3-butanoyl)-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(2,4-difluorophenyl)methyl-4-(3-butanoyl)-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 15 1-(4-chlorophenyl)methyl-4-(3-butanoyl)-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(diphenyl)methyl-4-(3-butanoyl)-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(*m*-trifluoromethyl)phenyl-4-(3-butanoyl)-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 20 1-methyl-4-(3-butanoyl)-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(1-methyl)ethyl-4-(3-butanoyl)-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 25 1-(2-methyl)propyl-4-(3-butanoyl)-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(2,4-difluorophenyl)methyl-4-(3-butanoyl)-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(4-chlorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 30 1-(diphenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(*m*-trifluoromethyl)phenyl-4-[(2-methyl)carboxyethyl]-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 35 1-methyl-4-[(2-methyl)carboxyethyl]-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(1-methyl)ethyl-4-[(2-methyl)carboxyethyl]-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.

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- 1-(2-methyl)propyl-4-[(2-methyl)carboxyethyl]-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(4-chlorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 5 1-(diphenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(*m*-trifluoromethyl)phenyl-4-[(2-methyl)carboxyethyl]-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-methyl-4-[(2-methyl)carboxyethyl]-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 10 1-(1-methyl)ethyl-4-[(2-methyl)carboxyethyl]-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(2-methyl)propyl-4-[(2-methyl)carboxyethyl]-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 15 1-(diphenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(6-aminoethyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(diphenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(6-aminoethyl-1-ynyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(diphenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(5-aminopentoxyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 20 1-(diphenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(5-guanidinopentyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(diphenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(4-guanidinobutoxy)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 25 1-(diphenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(5-aminopent-1-ynyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(diphenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(5-guanidinopent-1-ynyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(diphenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 30 1-(diphenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(diphenyl)methyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 35 1-(diphenyl)methyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(4-chlorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(6-aminoethyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.

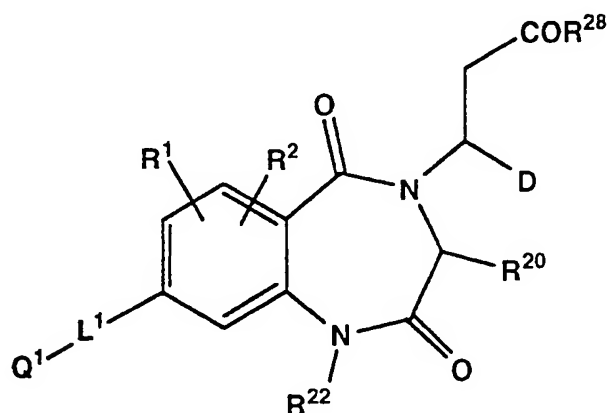
- 1-(4-chlorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(5-aminopentoxy)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(4-chlorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 5 1-(4-chlorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(4-chlorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(4-chlorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 10 1-methyl-4-[(2-methyl)carboxyethyl]-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-methyl-4-[(2-methyl)carboxyethyl]-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-methyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-methyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 15 1-(1-methyl)ethyl-4-[(2-methyl)carboxyethyl]-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(1-methyl)ethyl-4-[(2-methyl)carboxyethyl]-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(1-methyl)ethyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(1-methyl)ethyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 20 1-(2-methyl)ethyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(2-methyl)ethyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(2-methyl)propyl-4-[(2-methyl)carboxyethyl]-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(2-methyl)propyl-4-[(2-methyl)carboxyethyl]-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 30 1-(2-methyl)propyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(2-methyl)propyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(2,4-difluorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 35 1-(2,4-difluorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.

- 1-(2,4-difluorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(2,4-difluorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(5-aminopentoxyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 5 1-(2,4-difluorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(2,4-difluorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(2,4-difluorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 10 1-(2,4-difluorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(4-chlorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 15 1-(diphenyl)methyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(*m*-trifluoromethyl)phenyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-methyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 20 1-(1-methyl)ethyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(2-methyl)propyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 25 1-(2,4-difluorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(3-(butyryl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(3-(butyryl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(5-guanidinopentyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 30 1-(3-(4-aminobutyryl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(3-(4-guanidinobutyryl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(5-guanidinopentyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 35 1-(3-(6-aminohexoyl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(3-(6-guanidinohexoyl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(5-guanidinopentyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.

- 1-(3-(6-aminopentoyl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(3-(6-guanidinopentoyl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(5-guanidinopentyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 5 1-(3-(6-aminopropionyl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(3-(6-guanidinopropionyl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(5-guanidinopentyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(3-(6-aminoacetyl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 10 1-(3-(6-guanidinoacetyl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(5-guanidinopentyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(3-(4-guanidinobenzoyl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 15 1-(4-(6-aminopropionyl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(3-(4-aminobutyl)phenyl)methyl-4-(2-carboxyethyl)-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 1-(4-carboxyphenyl)methyl-4-(2-carboxyethyl)-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.
- 20

## 2. Nonpeptidyl $\alpha_v\beta_3$ inhibitors

- An alternative embodiment of the invention comprises a compound represented by formula VII capable of inhibiting binding of the osteoclast  $\alpha_v\beta_3$  receptor to its native *in vivo* ligands.
- 25



(VII)

where  $R^1$ ,  $R^2$ ,  $R^{20}$ ,  $R^{22}$ ,  $R^{28}$ ,  $Q^1$ , and  $L^1$  are defined above.

- 30 Compounds represented by formula VII are potent inhibitors of the vitronectin-vitronectin receptor (Vn-VnR) interaction, typically yielding  $IC_{50}$ 's in a Vn-VnR ELISA of

between 150 and 200 nM. These compounds show roughly a 20-fold selectivity when compared with the  $\text{GPII}_{\text{b}}\text{III}_{\text{a}}$ -Fg inhibition.

#### 5 D. Methods of Making

Compounds of the present invention can be prepared by many methods, employing standard chemical methodologies described and referenced in standard textbooks (e.g. March, J. "Advanced Organic Chemistry" McGraw-Hill, New York, 1977; Collman, J.P., Hegedus, L.S., Norton, J.R., Finke, R.G. "Principles and Applications of Organotransition Metal Chemistry" University Science, Mill Valley, 1987; Larock, R.C. "Comprehensive Organic Transformations" Verlag, New York, 1989). In the description that follows standard abbreviations as recommended by the *Journal of Organic Chemistry* (see "Guidelines for Authors" in any volume) are employed unless otherwise specified.

##### 1. The Q Group

15 The nitrogen containing substituents Q or a precursor thereof may added to the linker L and the combination Q-L-, usually in protected form, may be bonded to ring A. Alternatively, Q or Q plus a portion of L may be added to L or a portion thereof after ring A and the rest of the molecule have been formed. Q itself may be prepared and bonded to ring A by standard methods published in both the scientific and patent literature (see e.g. US patents; 20 4,992,542, 4,997,936, 4,194,047, 5,003,076, 5,063,207, 5,063,208, and 5,079,357, and references cited therein).

In the description immediately following, addition of Q-L- to the benzene moiety of 3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione is specified. It will be understood that these same procedures may be applied to other ring systems encompassed by formula I.

##### 25 2. The benzodiazepinedione

The key intermediate is the substituted 3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione 1. The benzodiazepinedione 1 ( $\text{R} \neq \text{H}$ ) may be prepared using a triply convergent approach from the isatoic anhydride 2, the b-alanine ester 3, and the a-haloacetyl halide 4 (Scheme 1). The benzodiazepinedione 1 ( $\text{R} = \text{H}$ ) may be prepared using a doubly convergent approach from the isatoic anhydride 2 ( $\text{R} = \text{H}$ ) and the substituted N-carboxymethyl b-alanine ester 5 (Scheme 2). 30

Methods for preparation of the isatoic anhydrides 2, b-alanine esters 3, and a-halo acetyl halides 4 (Scheme 1), are known in the art and a number of them are available from commercial sources such as Aldrich Chemical Co. A reaction sequence, similar to that shown in Scheme 1 to prepare 3,4-dihydro-1H-1,4-benzodiazepine-2,5-diones has been previously described (Lee, C.M. *J. Heterocyclic Chem.* 1: 235 (1964). Briefly, the isatoic anhydrides are converted to N-(2-aminobenzoyl)-b-alanine esters 6 by allowing 2 to react with 3, or its salt, in the presence an organic base. The isatoic anhydride 2 may be substituted or unsubstituted, but the 5-iodo isatoic anhydride is preferred (2,  $\text{R}^a = \text{I}$ ). The b-alanine ester 3 may also be substituted or unsubstituted as the free amine or, more conveniently, as its salt (e.g. HCl). For 35

example, the reaction may be carried out with b-alanine ethyl ester hydrochloride, ethyl 3-phenyl-3-aminopropionate p-tosylate, aspartyl-valine dibenzyl ester hydrochloride, or the like. Generally, the reaction is conducted in a dry polar aprotic solvent, such as dimethylformamide or the like, in the presence of an equimolar amount, or up to 30% excess, of a tertiary amine as the organic base, e.g. triethylamine. Alternatively, the reaction may be catalysed by dimethylaminopyridine (Venuti, M.C. *Synthesis* 266 (1982)). The reaction requires between about 0.5 to 4 hours at temperatures between about room temperature and 100°C. Preferably, the reaction is conducted at 50°C for about 2 hours in dimethylformamide as the solvent.

The products are isolated and purified by conventional methods, typically by solvent extraction into a compatible solvent. The product may be further purified by column chromatography or other appropriate methods.

If the anilino nitrogen in N-(2-aminobenzoyl)-b-alanine esters 6a is unsubstituted such that R=H, this position can be alkylated by allowing 6b (R=H) to react with a molar amount, but up to 50% excess, of an alkyl halide to produce 6 (R≠H). Generally, the reaction can be accomplished in a polar aprotic solvent, in the presence of a molar amount, or up to 50% excess, of an organic base. Preferably, the reaction will be conducted in dimethylformamide as solvent and 2,6-lutidine as the organic base. The reaction is allowed to proceed at temperatures between about room temperature to 200°C for about 0.5 hours to 2 days. Preferably, the reaction will be run at 80°C for 16 hours. The products are isolated and purified by conventional methods, typically by solvent extraction into a compatible solvent. The product may be further purified by column chromatography or other appropriate methods.

Alternatively, 6 (R≠H) may be prepared by allowing 6a to react with an aldehyde or ketone, under dehydrating conditions, in the presence of a catalytic amount of acid followed by reduction with a trialkylsilane and a strong acid. Generally, the first of these reactions can be accomplished in a non-polar aprotic solvent, such as toluene, and an organic acid, such as p-toluenesulphonic acid at temperatures between 80-160°C for times between 1 to 60 mins. Preferably, the reaction will be conducted at reflux for 10 mins. The product may be isolated by crystallization or carried on to reduction without further purification. Reduction can be achieved by the addition of an excess of trialkylsilane, such as triethylsilane, in the presence of an excess of a strong acid, such as trifluoroacetic acid, at temperatures between -30 to 30°C for times between 0.5 to 48 hours with a non-reactive organic solvent, such as dichloromethane. Preferably, reduction is run at 0°C for 2 hours with dichloromethane as a co-solvent. The products are isolated and purified by conventional methods, typically by crystallization from suitable solvents. The product may be otherwise purified by solvent extraction into a compatible solvent followed by column chromatography or other appropriate methods.

Conversion of N-(2-aminobenzoyl)-b-alanine esters 6 to the benzodiazepinedione 1 (R≠H) involves first the acylation of the anilino nitrogen, followed by a base promoted ring formation. The acylation reaction can be achieved by either reaction of 6 with an α-haloacyl halide in the presence of an equimolar amount, or up to 30% excess, of an organic base, such as

triethylamine, or with an  $\alpha$ -haloacid in the presence of an amide coupling reagent, such as dicyclohexylcarbodiimide. Acylation of 6 with an  $\alpha$ -haloacyl halide may also be conducted in the absence of an organic base and can be run with water in a biphasic reaction system. Generally, these reactions are conducted in a pre-dried, non-reactive organic solvent, such as

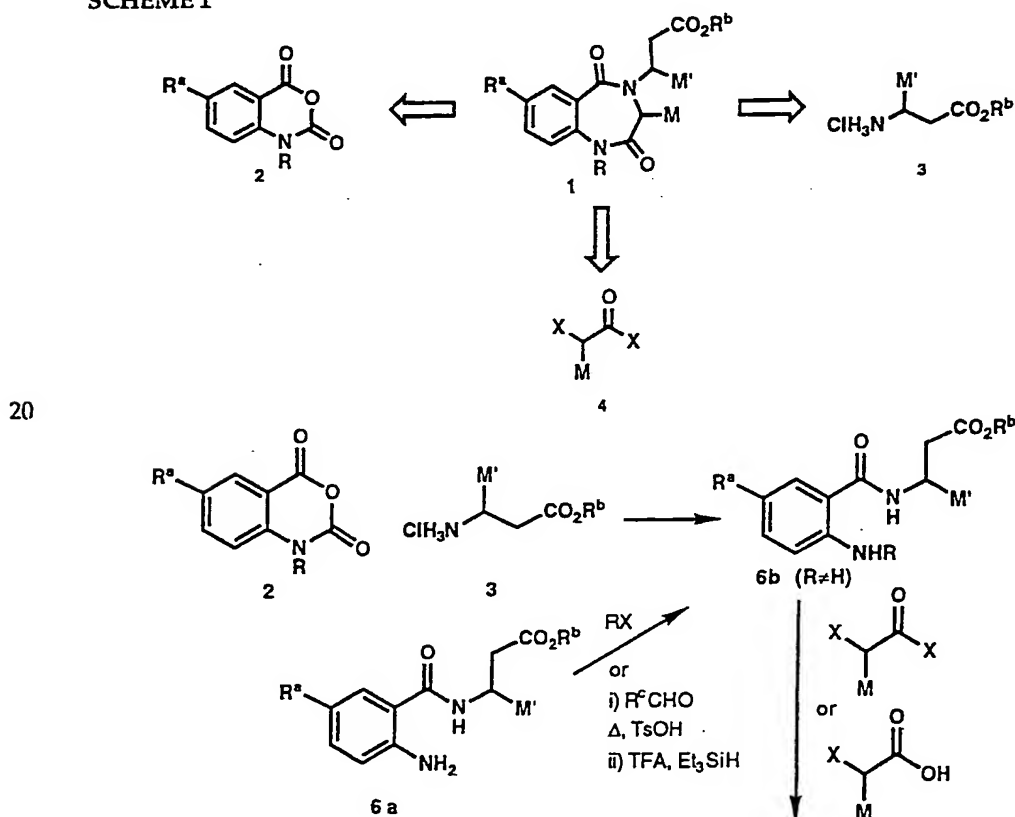
5 methylene chloride, diethyl ether, or tetrahydrofuran under a dry, inert atmosphere, such as nitrogen. Preferably, acylation with an  $\alpha$ -haloacylhalide will be run in methylene chloride, whereas acylation with an  $\alpha$ -halo-acid will be run in tetrahydrofuran. The reactions are allowed to run at temperatures between about 0°C to room temperature for times between about 0.5 to 24 hours. Preferably, the reactions will be allowed to run for 2 hours at room

10 temperature. The products are isolated and purified by conventional methods, typically by solvent extraction into a compatible solvent. Generally, further purification of the acylated products 7 is not required.

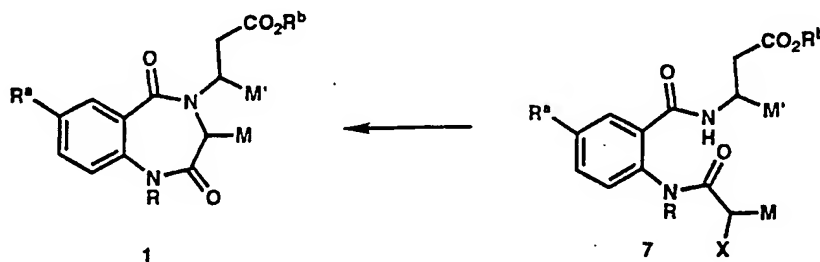
Cyclization of the acylated N-(2-acylaminobenzoyl)-b-alanine esters 7 may be achieved by reaction with a base such as an alkali metal alkoxide, hydride, or carbonate in a polar

15 solvent at temperatures between about 0 and 100°C for about 0.5 to 2 hours. For example, a solution of the N-(2-aminobenzoyl)-b-alanine ester 7 is commonly added over a period of approximately 15 to 60 minutes to a slurry of an alkali metal hydride in an appropriate solvent, cooled to 0°C.

SCHEME 1





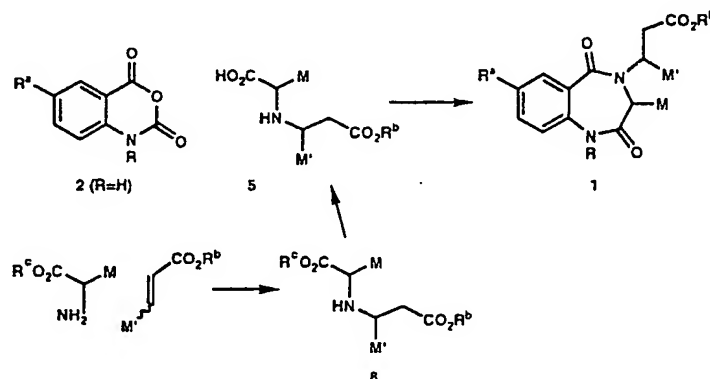


It is preferable that the solvent will be a polar aprotic solvent such as dimethylformamide and the alkali metal hydride be sodium hydride or calcium carbonate. Once the addition is complete, the reaction mixture is generally allowed to warm to room temperature and run for an additional 60 to 105 minutes after which the reaction is neutralized by the addition of a solution of an acid, such as 10% citric acid or the like, and the solvent evaporated. The product is then isolated by solvent extraction and further purified by column chromatography.

In Scheme 2, the substituted N-carboxymethyl b-alanine esters 5 are also known in the art and may be prepared in two steps by first reaction of an acrylate ester and a-amino acid ester to give 8, followed by selective conversion of the glycine ester group to its corresponding acid. For example, allowing glycine benzyl ester to react with ethyl acrylate in an alcoholic solvent such as methanol for about 1 day yields N-(carboxymethyl benzyl ester)-b-alanine ethyl ester. Removal of the benzyl ester by hydrogenation in the presence of a catalytic amount 10% palladium on carbon provides N-carboxymethyl-b-alanine ester 5 (M'=H, R<sup>b</sup>=ethyl).

Preparation of the benzodiazepinedione 1 (R=H) may be accomplished by allowing 5 to react with 2 (R=H). The reaction can be conducted in either a

**SCHEME 2**



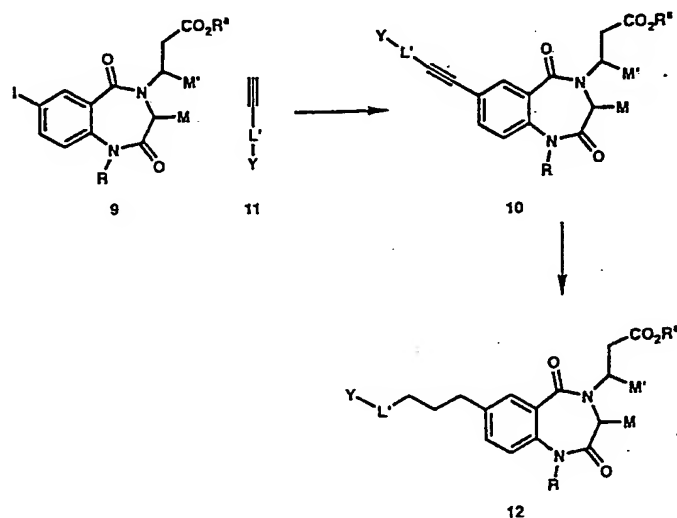
polar aprotic solvent, such as dimethyl sulfoxide, or weak organic base, such as pyridine, at temperatures of 100-200°C for about 2-86 hours. Preferably, the reaction will be run in pyridine at refluxing temperature for about 24 hours. The product is then isolated by dilution with an appropriate organic solvent, such as ethyl acetate, and washed with aqueous acid, such as 10% citric acid, and aqueous base, such as saturated sodium bicarbonate. The products can be further purified by column chromatography or crystallization from an appropriate organic solvent to provide the compounds 1 (R=H).

Reaction Scheme 3 outlines the preparation of the compounds of the present invention from compound 9 [ $R^a=I$ ].

Conversion of the compounds having the structural formula 9 to compounds having the structural formula 10, provides those compounds classified in Groups 1 and 2. This may be accomplished by allowing 9 to react with an alkyne 11 in the presence of a palladium (II) salt, a copper (I) salt, and an organic base. Preferably, the alkyne 11 is substituted in a way such that the Y group can be synthetically converted to the positively charged Q group of structural Formula I. More preferably, the Y group is a protected form of the positively charged Q group of structural Formula I. For example, the alkyne may be a N-Boc-amino alkyne, a benzo or alkynitrile alkyne, a nitrobenzo alkyne, or the like. L' is L less the functional groups shown below in scheme 3.

Generally, the reaction is allowed to run in a dry organic polar aprotic solvent, such as ethyl acetate or the like, to the exclusion of oxygen at temperatures between about

SCHEME 3

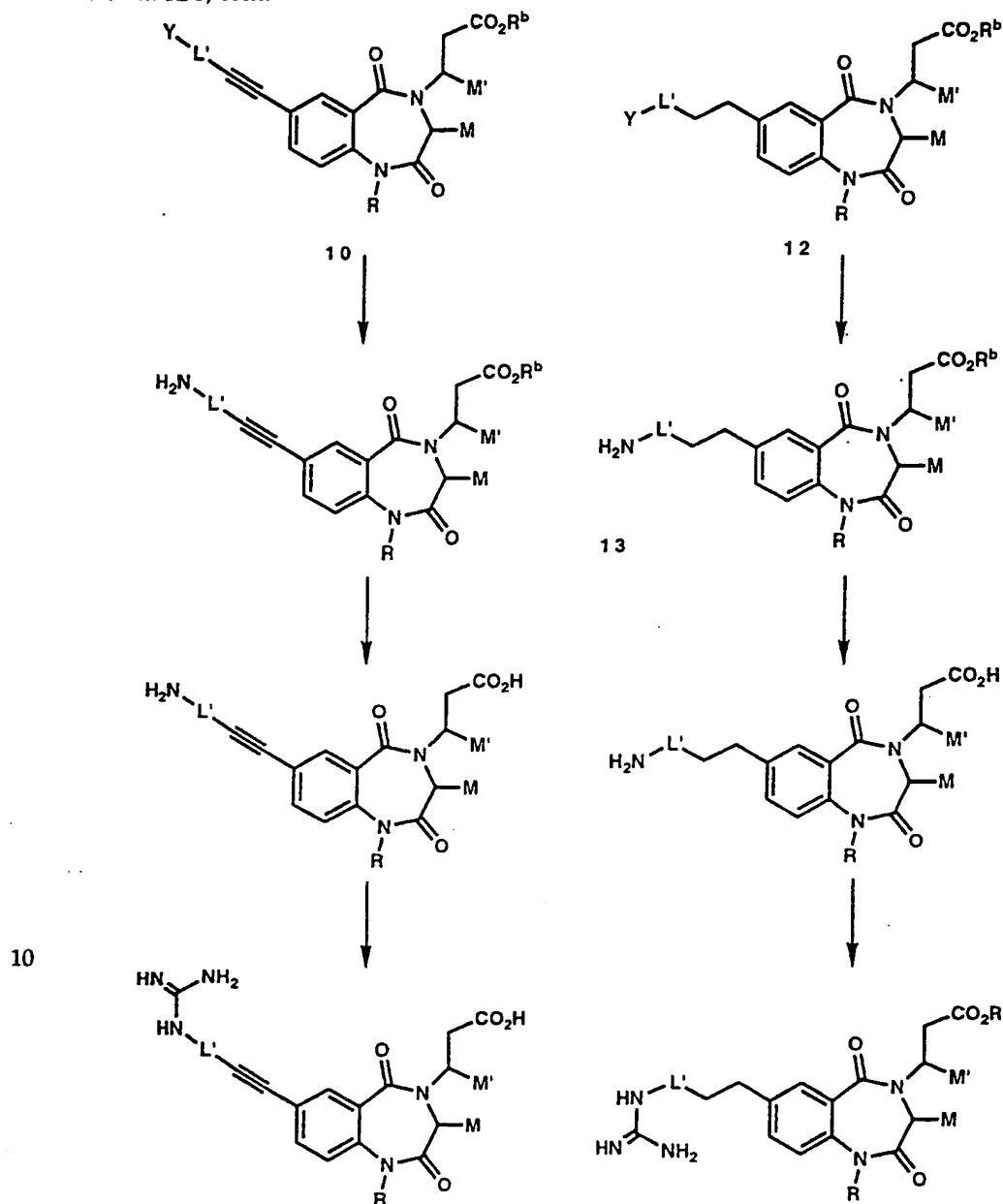


room temperature and 180°C for times between about 2 and 48 hours. Preferably, the reaction will run with a catalytic amount of palladium(II) salt at 10 molar percent and cupric(I) salt at 5 molar percent. A 2 fold excess of the alkyne 11 and a 5 fold excess of a tertiary amine as the organic base, such as triethylamine or the like, is preferred. For example, a mixture of the iodo arene 9, a 2 fold excess of the alkyne 11, 10 molar percent of bis(triphenylphosphine)palladium dichloride, 5 molar percent cupric iodide, 5 fold excess of triethylamine, and ethyl acetate under a dry, inert atmosphere, such as nitrogen, is allowed run for about 2 hours. The product can be isolated by solvent extraction in to a suitable organic solvent, such as ethyl acetate, and washed with a solution of 10% ethylenediaminetetraacetic acid and the solvent evaporated. The products may be further purified by column chromatography.

Preparation of the compounds classified in Groups 3 and 4 may be achieved by reduction of the alkyne moiety in compounds of the general formula 10 to yield the saturated compounds 12 by allowing the alkyne to stir under an atmosphere of hydrogen in the presence

- of a small amount of palladium on carbon. Typically, the reaction is run in an inert solvent, such as ethyl acetate, with a 5-10 molar percent by weight of 10% palladium on carbon at temperatures between about room temperature and 50°C for times between about 15 - 240 minutes. Preferably, the reaction is carried out at room temperature for 1 hour. The products
- 5 are isolated by filtration of the mixture through a filter agent, such as Celite®, and evaporation of solvent.

SCHEME 3, cont.



Conversion of the compounds of the general formula 10 or 12 in which Y is a N-(*tert*-butoxycarbonyl) amino protected moiety (N-BOC) to the amino esters 13 can be accomplished

by allowing the material to react with a strong acid. Generally, the reaction is carried out by mixing the N-BOC amino ester with a large excess of concentrated solution of a strong acid, such as hydrogen chloride, dissolved in an appropriate inert solvent, such as ethyl acetate. The reaction can be conducted at a temperature between about minus 30°C and room temperature for about 0.5 to 24 hours. The reaction can be run in the presence of a molar amount of trialkylsilane, such as triethylsilane or the like. Preferably, the reaction will be carried out at room temperature for 2 hours with a molar equivalent of triethylsilane in ethyl acetate. The solvent is evaporated and the products can be further purified by high-pressure liquid chromatography using a reverse phase column.

Conversion of the compounds of the general formula 10 or 12 in which the Y functional group is an aryl nitrile to the amino esters 13 can be achieved by allowing the aryl nitriles 10 or 12 to react with dicobalt octacarbonyl in the presence of trimethylsilane. The reduction of a benzonitrile to a benzylamine is a known reaction (Murai, T.; Sakane, T.; Kato, S. *Tetrahedron Lett.* 26: 5145-5148[1985]) and is generally carried out in an inert solvent, such as toluene or the like, at 60°C for 20 hours with an 8 molar percent of dicobalt octacarbonyl and a 10 fold excess of trimethylsilane. The solvent is evaporated and the resulting material diluted with methanol and allowed to react with a 5 fold excess of potassium fluoride. The products are then isolated by solvent extraction and further purified high-pressure liquid chromatography using a reverse phase column.

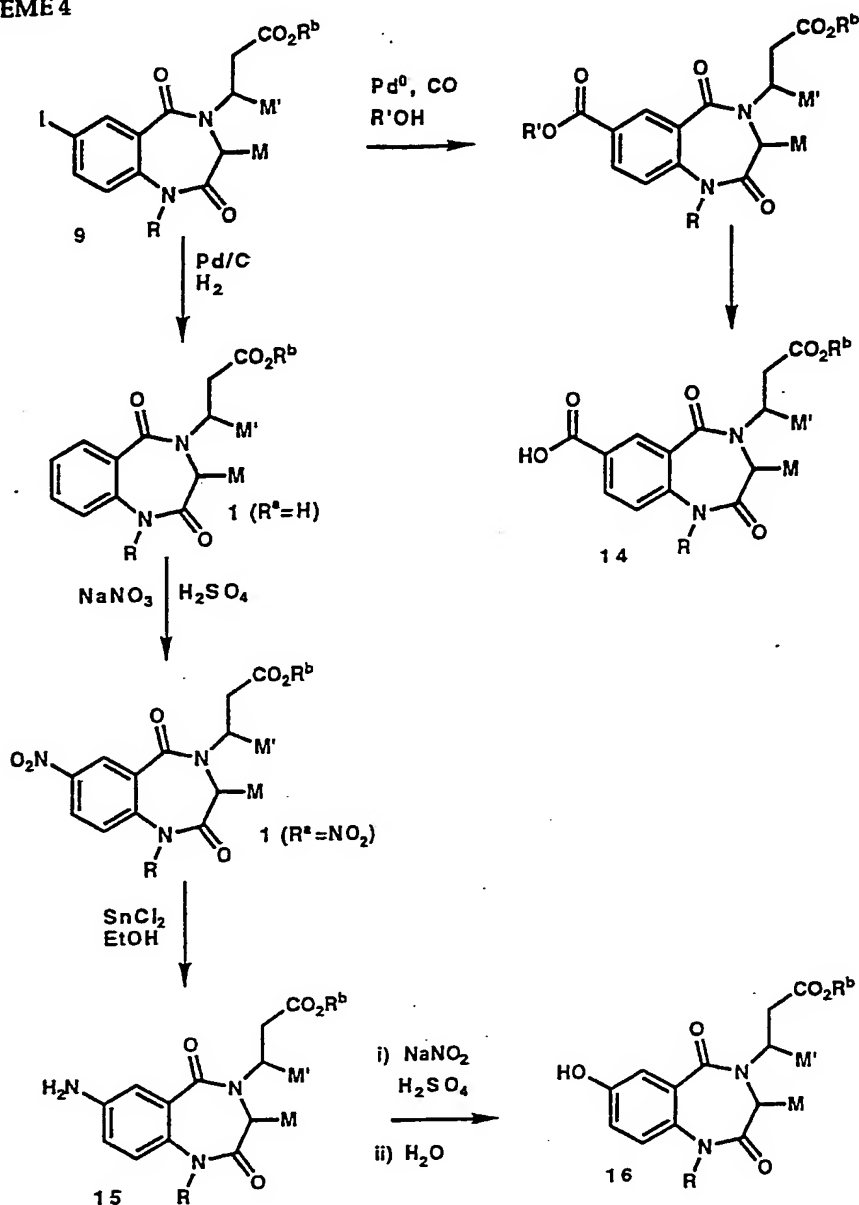
Conversion of the compounds of the general formula 10 or 12 in which the Y functional group is a nitroarene, to the amino ester 13 can be executed by a selective reduction (Bellamy, F.D.; Ou, K. *Tetrahedron Lett.* 25: 839-842[1984]). The reaction is generally run with a five molar excess of stannous chloride dihydrate in either ethyl acetate or ethanol as the solvent at temperatures between about 50 and 100°C for times between about 15-120 minutes under an inert atmosphere, such as nitrogen. Preferably, the reaction is carried out at 70°C in ethanol for approximately 30 minutes. The products are then isolated by solvent extraction and further purified by high-pressure liquid chromatography using a reverse phase column.

Conversion of the amino esters to their corresponding amino acids of the structural formula I involves saponification using well known conditions and reagents. For example, an aqueous solution of a strong alkali metal base, such as sodium hydroxide, lithium hydroxide or the like, is added to an alcoholic solution of the ester. Alcohols which may be used as the solvent for this reaction may include, for example, methanol, ethanol, and isopropanol, but, methanol is preferred. The preferred base is sodium hydroxide at a concentration between about 1 to 6N, though 2N is preferred. The reaction may be conducted at a temperature between about 0 to 50°C for times between about 10 to 60 minutes. Preferably, the reaction is carried out at room temperature for 30 minutes after which the reaction is neutralized with a concentrated solution of a strong acid, such as hydrochloric acid or the like, and the solvent evaporated. The products are isolated by high-pressure liquid chromatography using a reverse phase column.

The conversion of the amino acids to their corresponding guanidino acids of the structural formula I is a known reaction (Kim, K. Lin, Y.-L.; Mosher, H.S. *Tetrahedron Lett.* 3183-3186 [1988]). The reaction can be accomplished by allowing the amino acid to react with aminoiminomethanesulfonic acid. Generally, the reaction can be conducted with a equimolar to a 10 fold molar excess of aminoiminomethanesulfonic acid at temperatures between 0 to 50°C for times between about 15 to 120 minutes in a polar protic solvent, such as methanol, water or the like. The solution, prior to the addition of aminoiminomethanesulfonic acid, may be made neutral or basic by the addition of weak base, such as an alkali metal carbonate. Preferably, for alkyl amines the amino acid will be allowed to react with a 5 fold excess aminoiminomethanesulfonic acid at room temperature for 30 minutes with 5% potassium bicarbonate in water as the reaction medium, whereas for aryl amines the amino acid will be allowed to react with an equimolar amount of aminoiminomethanesulfonic acid at room temperature for 1 hour in methanol. Generally, the reaction mixture will be made acidic by the addition of a dilute solution of an acid, such as acetic acid and the solvent evaporated. The products are isolated by high-pressure liquid chromatography using a reverse phase column.

Preparation of the compounds classified as Groups 4-11 may also be prepared from the key intermediate 9. Conversion of the iodoarene 9 to a synthetic precursors of those compounds found in Groups 5-11 namely compounds 15-18, may be accomplished by a multistep reaction sequence (Scheme 4). First, the iodoarene is carbonylated in an alcoholic solvent (R'OH) under a carbon monoxide atmosphere utilizing a palladium (0) catalyst. The alcoholic solvent (R'OH) must be chosen so as to allow for the selective removal of the R' group of the diester 14. For example, when R' = CH<sub>3</sub> and R<sup>b</sup> = *tert*-butyl, the alcoholic solvent is methanol, and this methyl ester functionality in 14 can be removed by mild basic hydrolysis to afford acid 15. The treatment of 15 with a molar equivalent of diphenylphosphoryl azide and a molar equivalent of a tertiary amine such as triethylamine at room temperature in a non-protic solvent such as dichloromethane affords the isocyanate 16. This may be hydrolyzed to the amine 17 by the addition of water to a solution of 16 in tetrahydrofuran at room temperature. Finally, conversion of the aniline 17 to its corresponding diazonium salt by treatment with nitrous acid and subsequent heating of the diazonium salt in the presence of water produces the phenol 18.

SCHEME 4

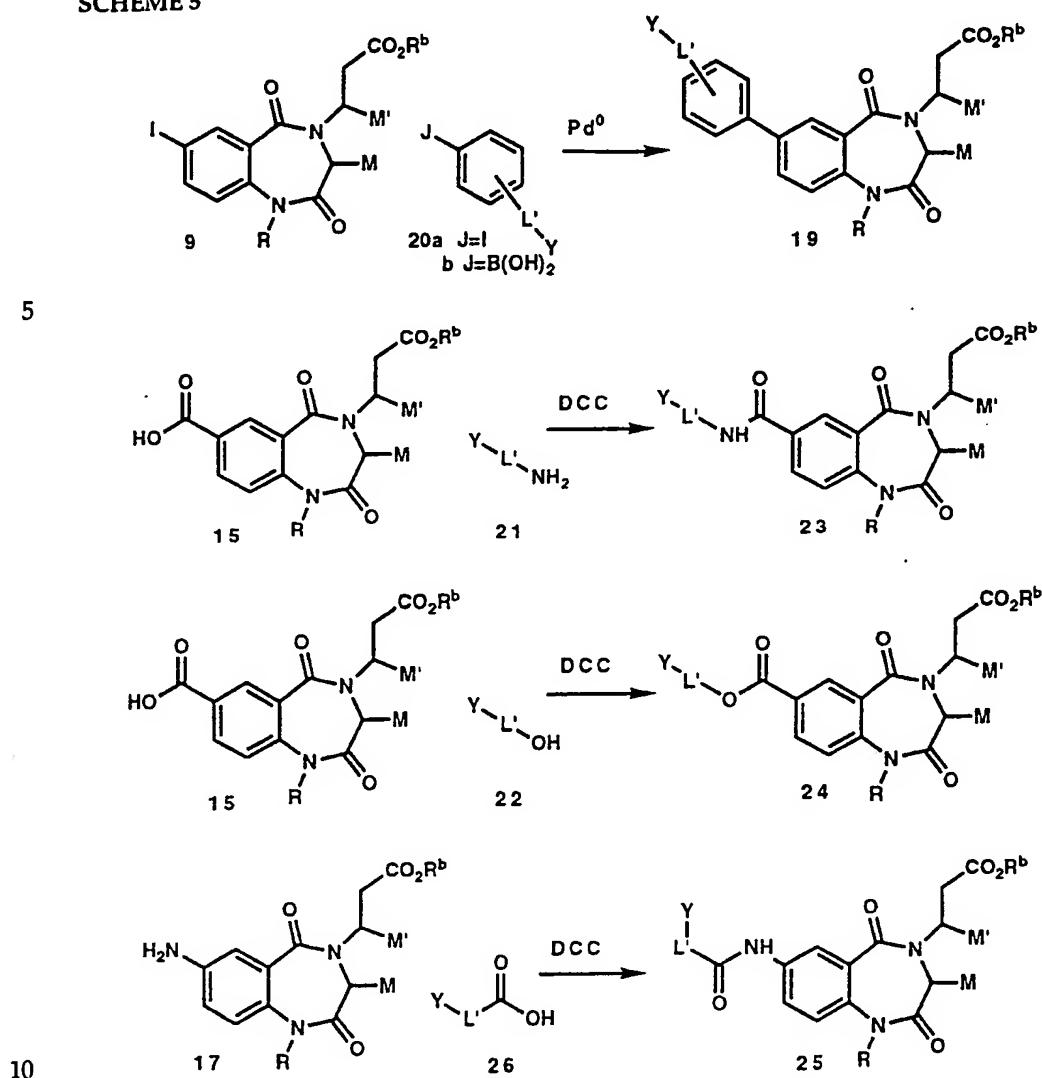


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Schemes 5 and 6 depict methods that may be used to introduce the linking groups of compounds classified as Groups 4-11 onto the benzodiazepinedione nucleus. For example, preparation of the biaryl adducts 19, precursors of Group 5 type compounds, may be accomplished by allowing 9 to react with substituted aryl molecules 20, where Y and L' are defined as above, in the presence of palladium(0). Allowing the carboxylic acid 15 to react with an amine 21 or alcohol 22 in the presence of a dehydrating reagent, such as dicyclohexylcarbodiimide (DCC) or the like, will yield the amide 23 and ester 24, forerunners of Group 10 and 11 type compounds, respectively. Similarly, 25 can be prepared from the aniline

17 by reaction with a carboxylic acid 26 in the presence of a coupling reagent, such as DCC or the like. Compounds of the general formula 25 are precursors of Group 5 type compounds.

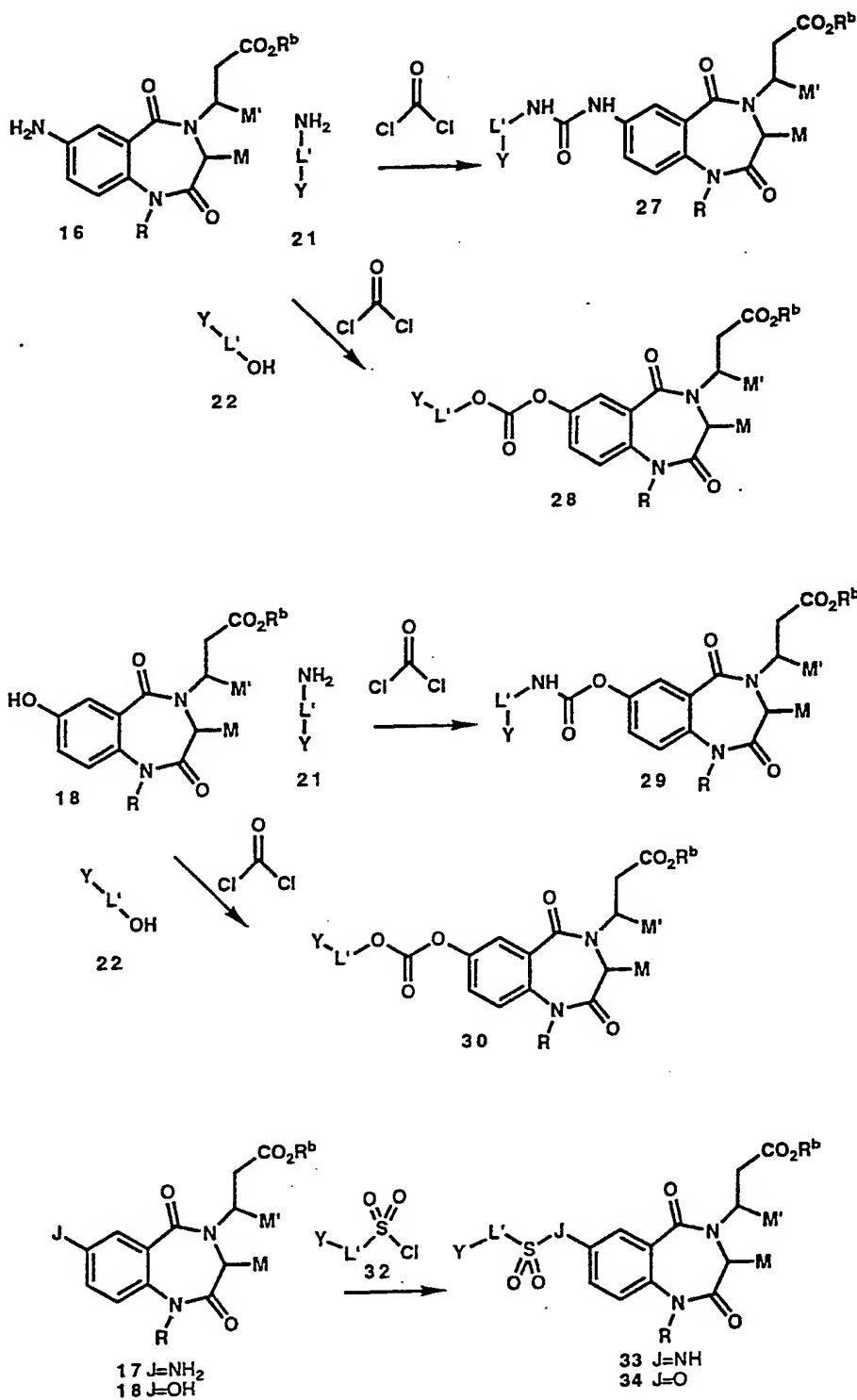
SCHEME 5



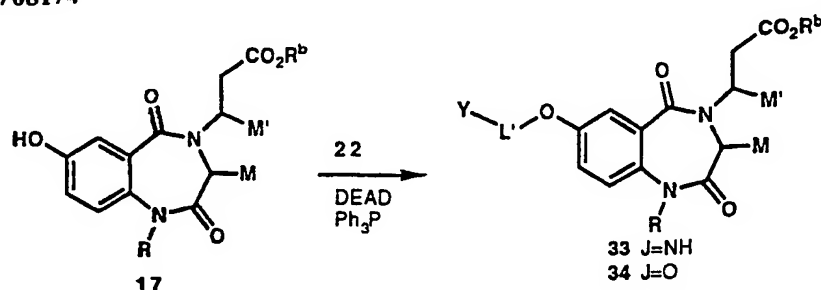
Conversion of aniline 17 into its corresponding urea 27 or carbamate 28 (Scheme 6) may be accomplished by reaction of the isocyanate 16 in the presence of an amine 21 or alcohol 22, respectively. Treatment of the phenol 18 with phosgene and an amine 21 or alcohol 22 yields the carbamate 29 and carbonate 30, respectively. Reaction of a sulfonylhalide 32 with either the aniline 17 or phenol 18 would produce the corresponding sulfonamide 33 and sufinyl ester 34, respectively. Allowing the phenol 18 to react with an alcohol 22 in the presence of diethylazodicarboxylate and triphenylphosphine will furnish 31, a precursor to Group 12 type compounds. The product 27 is a synthetic precursor to those compounds classified as Group 7, product 28 is a forerunner to Group 6 type compounds, product 29 will give rise to Group 8

type compounds, whereas, product 30 is a precursor to Group 9 type compounds. Products 33 and 34 are precursors to Group 15 and 16, respectively.

SCHEME 6



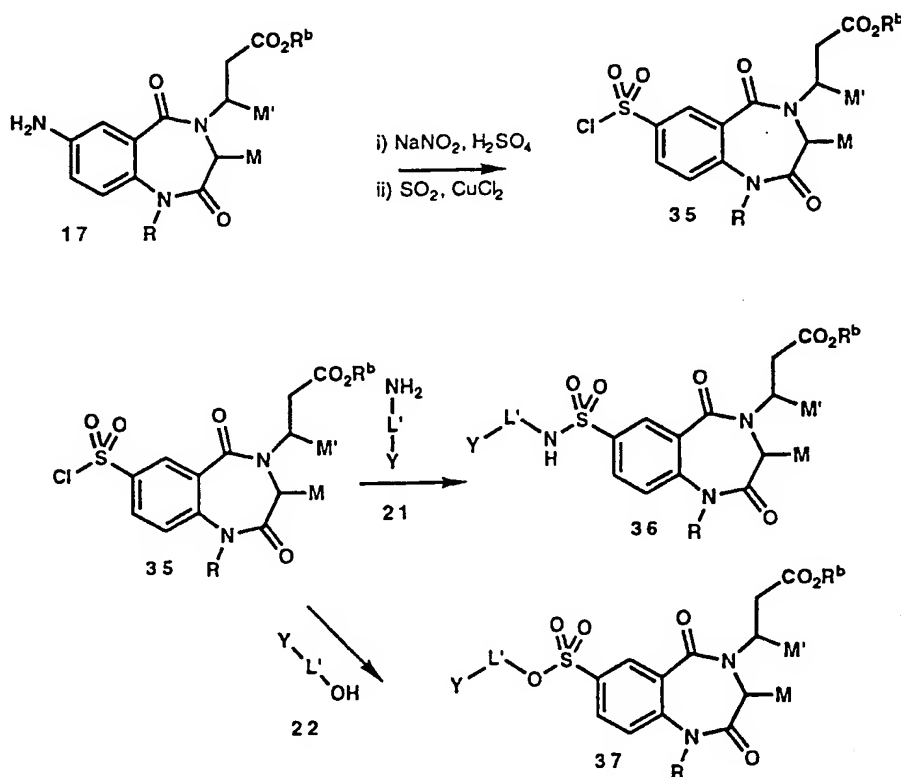


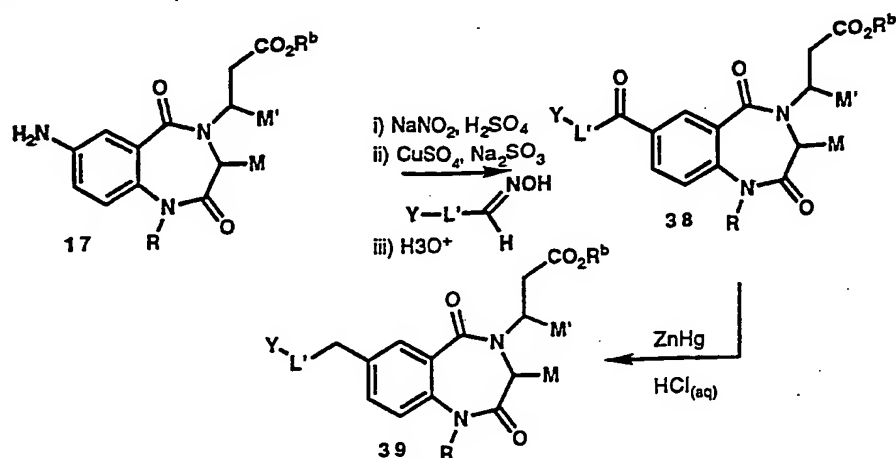


Scheme 7 depict methods that may be used in the preparation of the compounds classified as Groups 13, 14, 17 and 18. Diazotization of the aniline 17, by the reaction of 17 with nitrous acid, followed by treatment with sulfur dioxide in the presence of cupric chloride (Gilbert, *Synthesis*, 1969, 1-10) yields the sulfonyl chloride 35. Reaction of 35 with either the amine 21 or the alcohol 22 will provide the synthetic precursor to Groups 13 and 14, respectively, namely, the suphonamide 36 and sulfinyl ester 37. The diazonium salt prepared by diazotization of 17 when allowed to react with oximes in the presence of copper sulphate and sodium sulfite will give the corresponding oxime 38, which can be readily hydrolyzed to yield the aryl ketone 39 (Beech, *J. Chem. Soc.*, 1954, 1297). Clemmensen reduction of the ketone 39 using zinc amalgam and aqueous hydrogen chloride gives rise to the benzyl adduct 40 (Vedejs, *Org. React.* 1975, 22, 401). The products 39 and 40 are precursors to the compounds classified as Groups 17 and 18, respectively.

## SCHEME 7

15



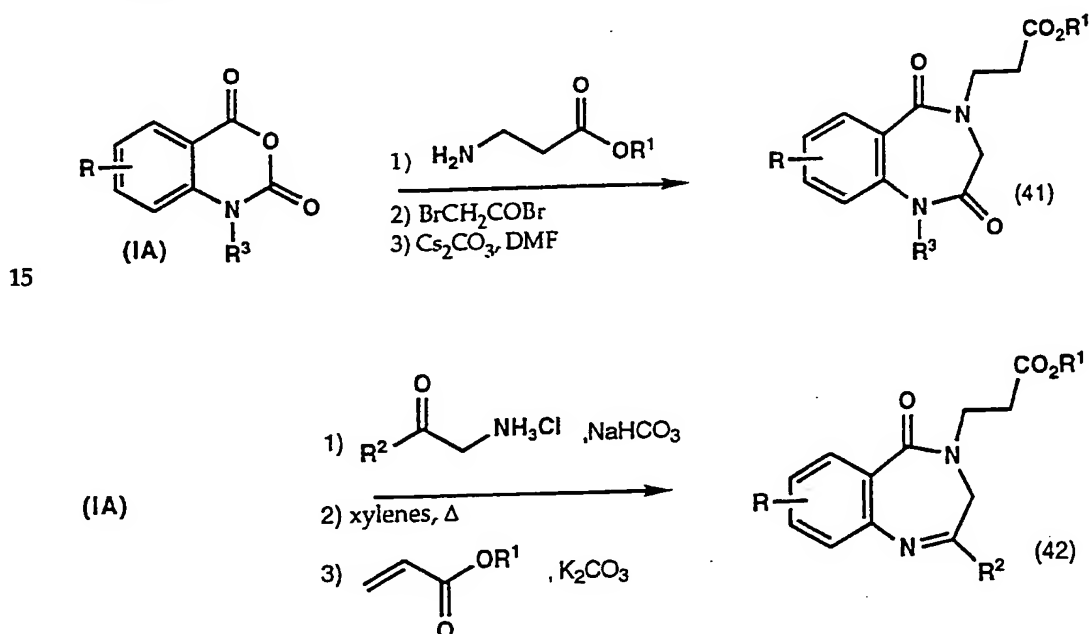


Each of the protected precursors 19, 23, 24, 25, 27, 28, 29, 30, 31, 33, 34, 36, 37, 39, and 40 may be treated in a similar way as described above for compounds 10 and 12 (Scheme 3) to prepare the compounds encompassed by this invention.

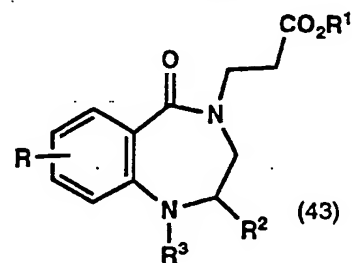
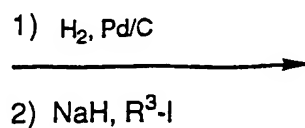
### 3. Other "6-7" Fused Ring Systems

Other "6-7" fused ring systems where each T, U, or G of formula I is independently selected from NH, N, NR, C=X, CR<sub>2</sub>, CHR, CH<sub>2</sub>, CR, CH, O, S, SO, or SO<sub>2</sub> can be prepared by methods known in the art. To provide guidance, however, specific representative synthetic routes are provided in scheme 8. It will be understood that other synthetic routes may be preferred when specific substituents or T-U-G are targeted.

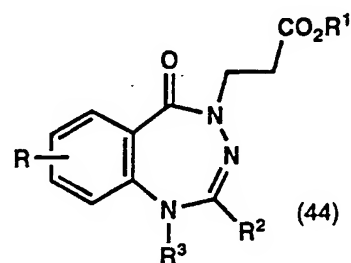
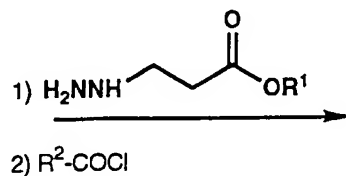
SCHEME 8



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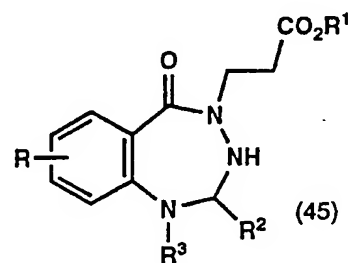
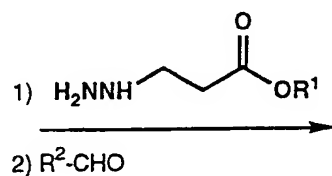


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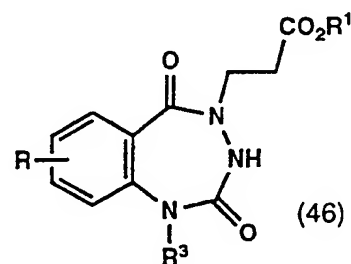
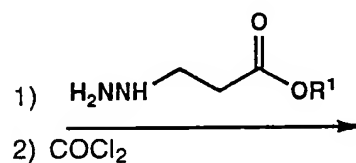


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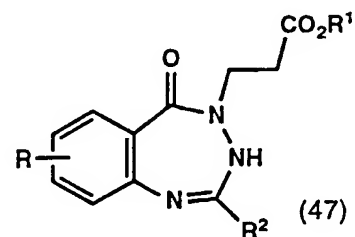
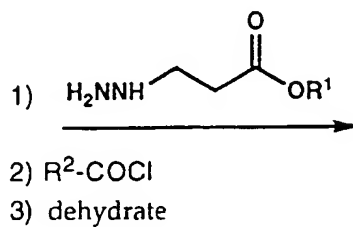


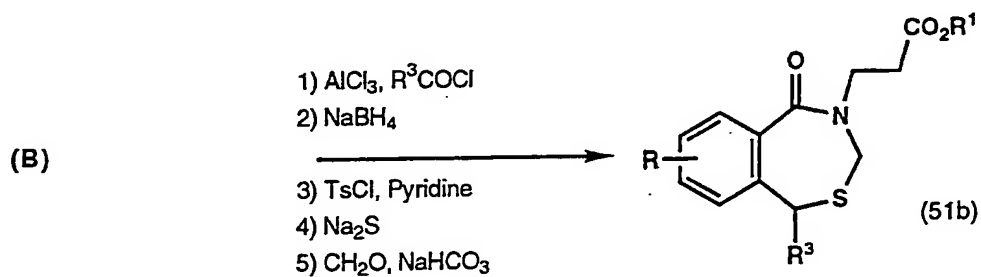
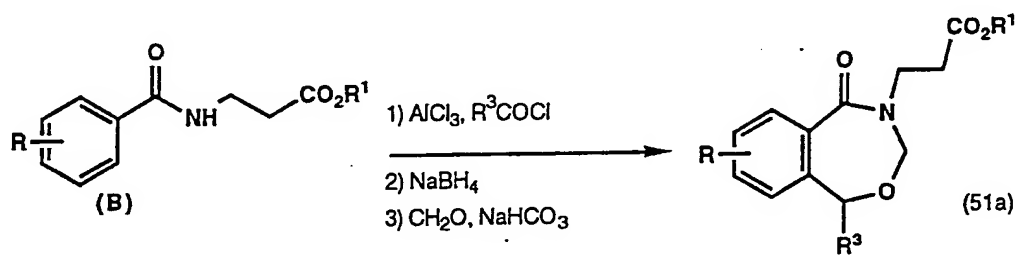
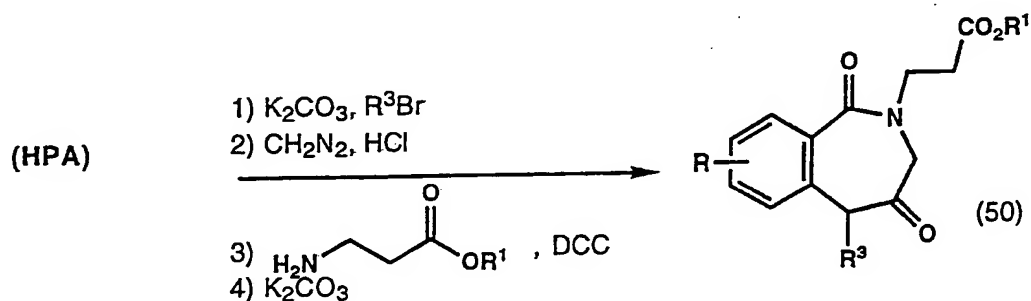
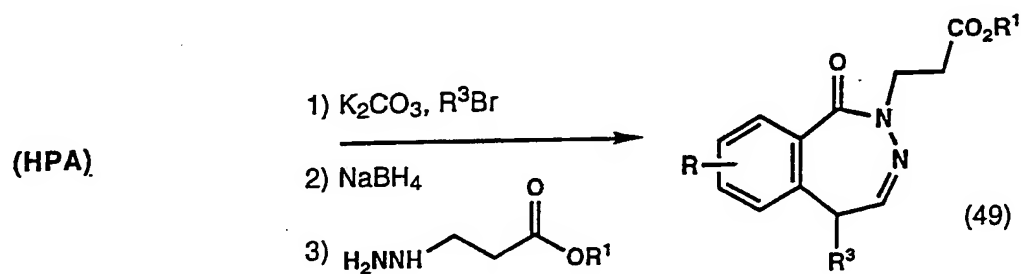
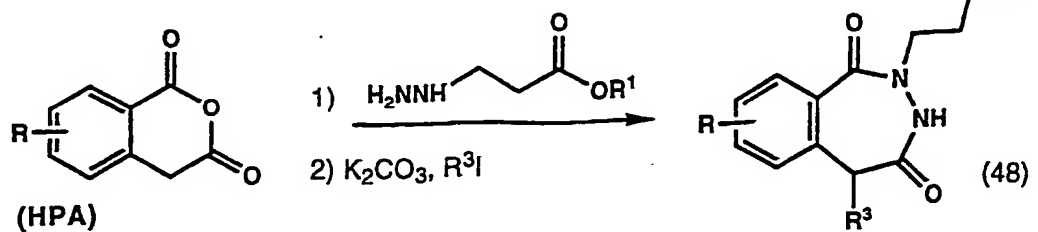
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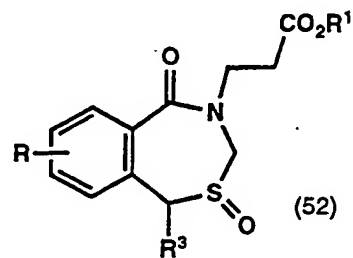
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(51b)

m-CPBA

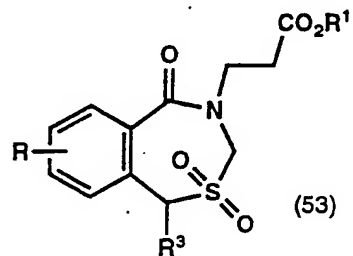


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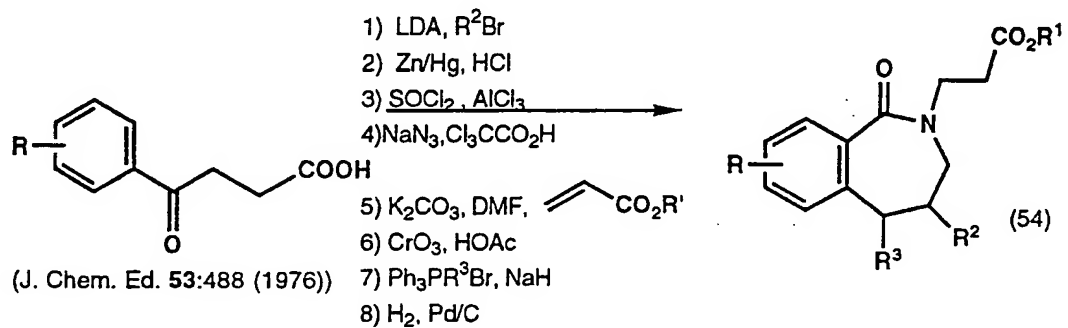
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(51b)

m-CPBA



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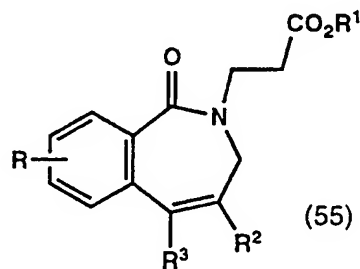


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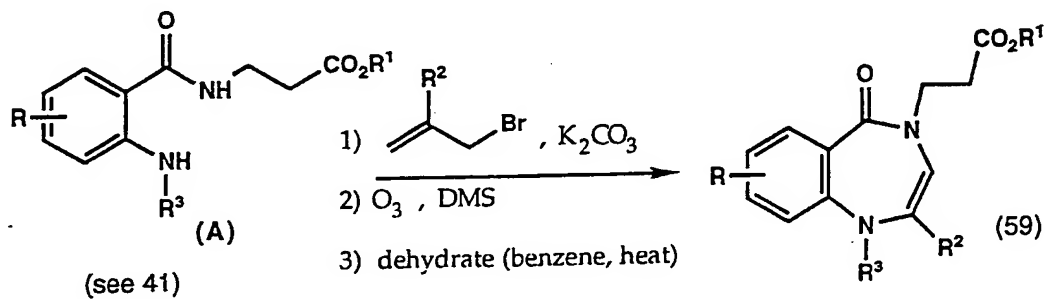
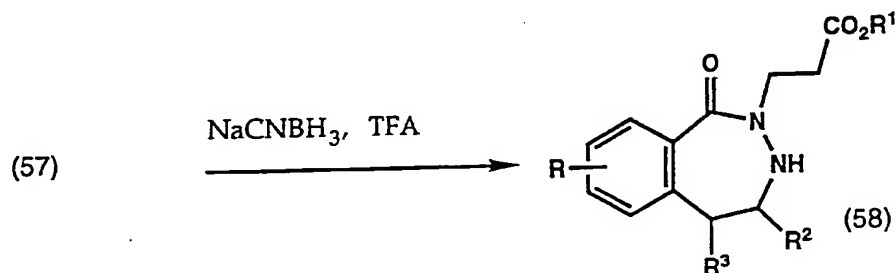
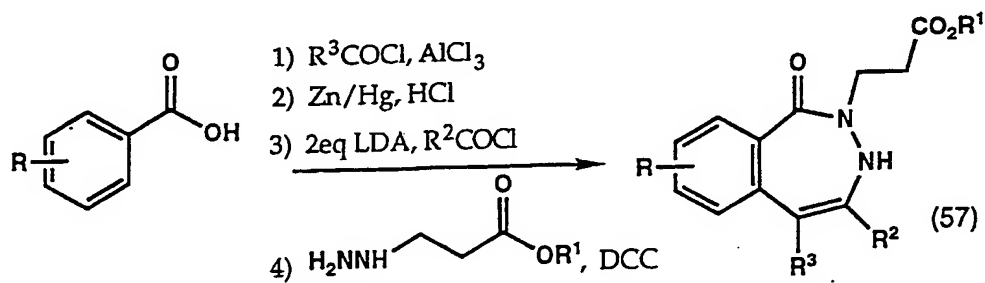
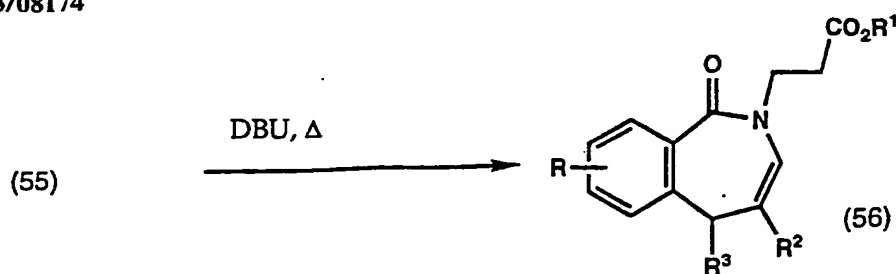
(54)

1) NBS, CCl<sub>4</sub>,  
benzoyl peroxide  
2) DBU

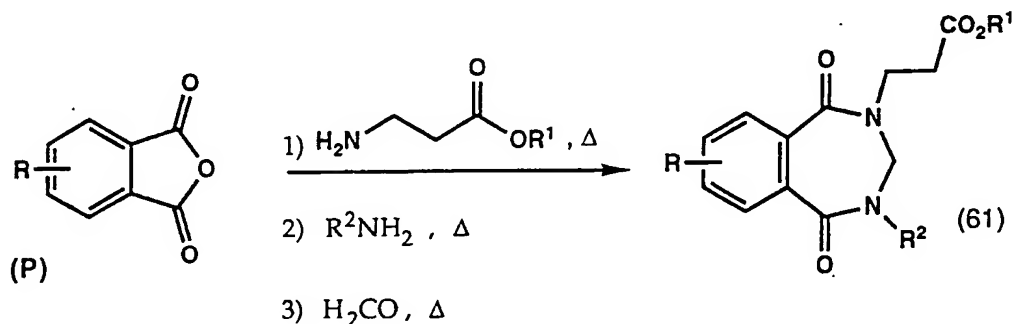
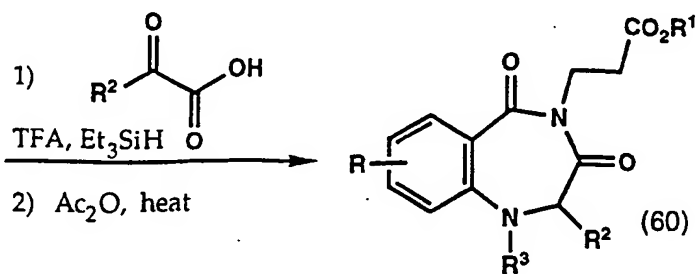


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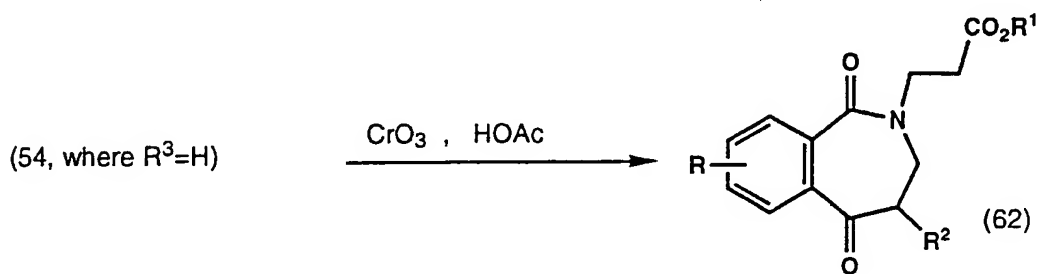
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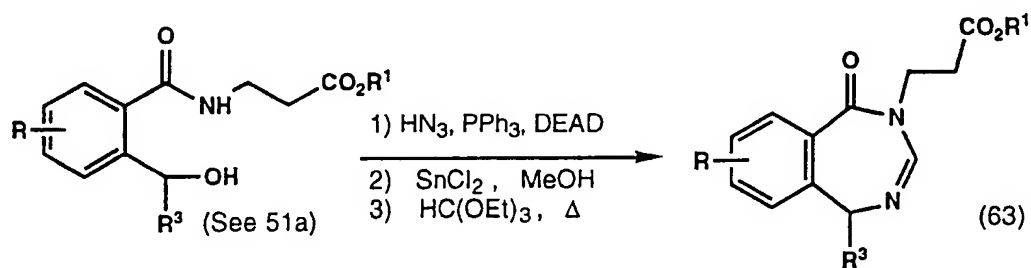
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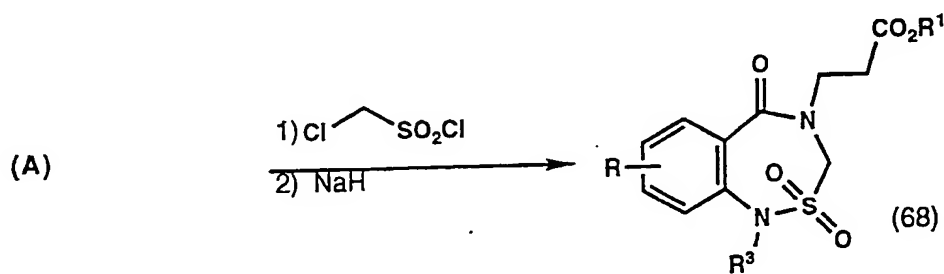
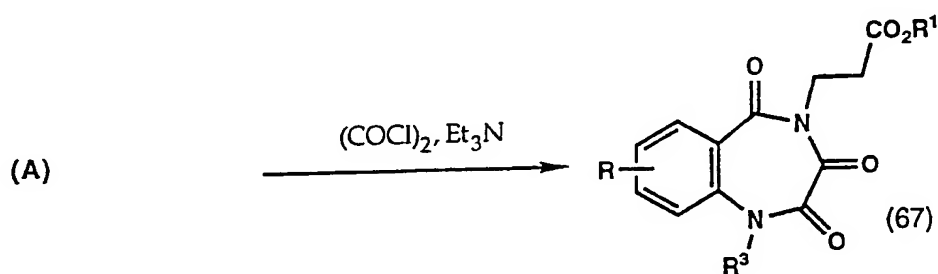
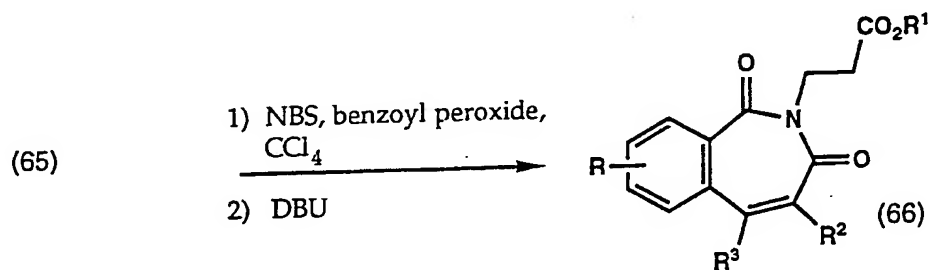
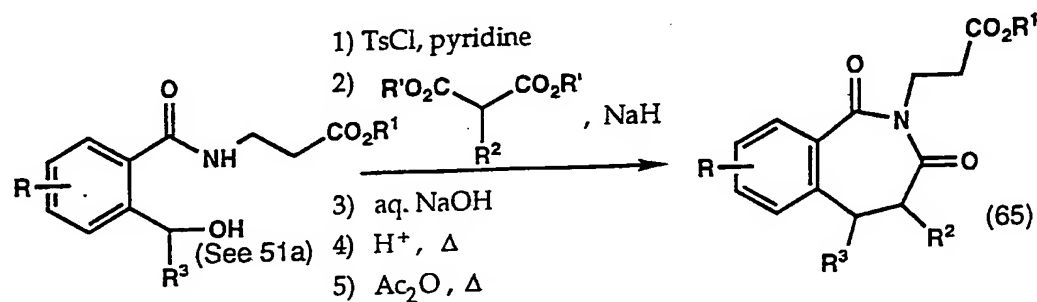
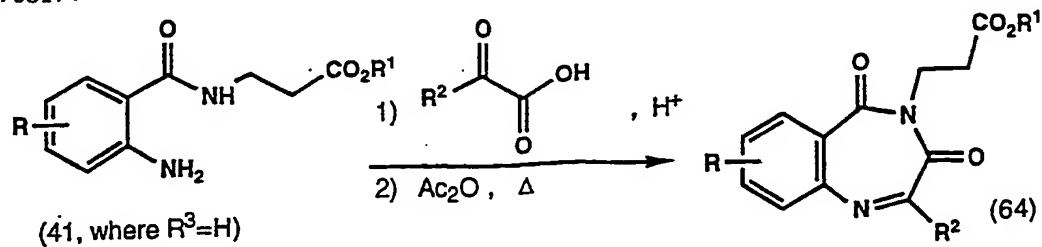
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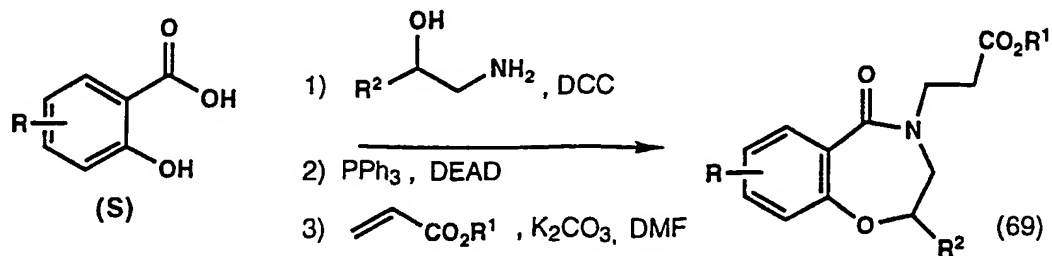
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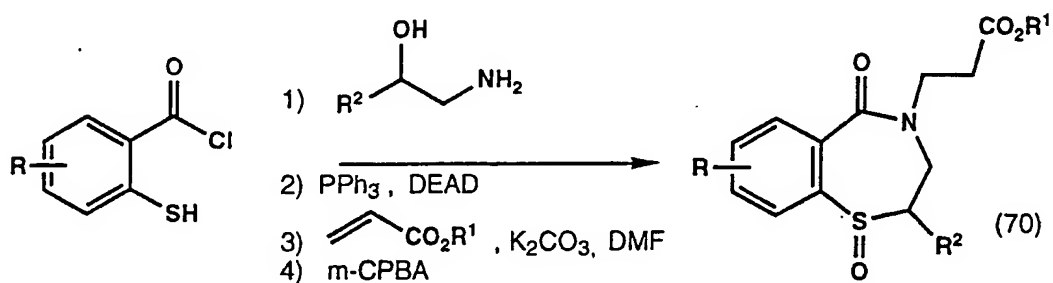
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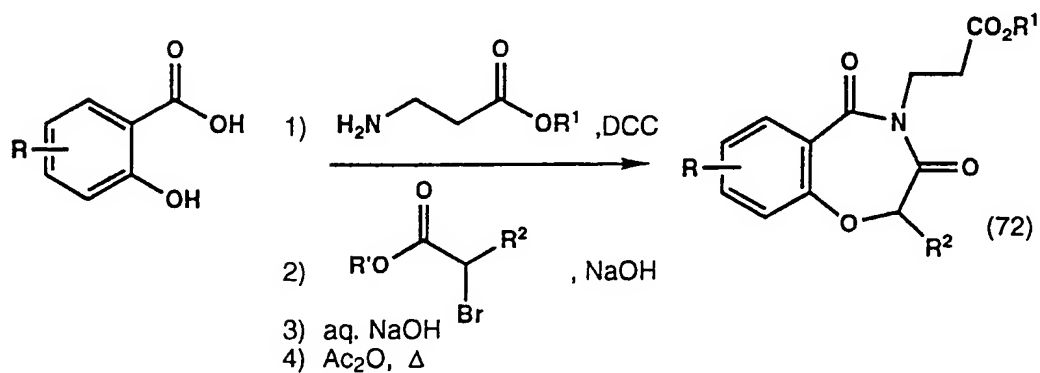
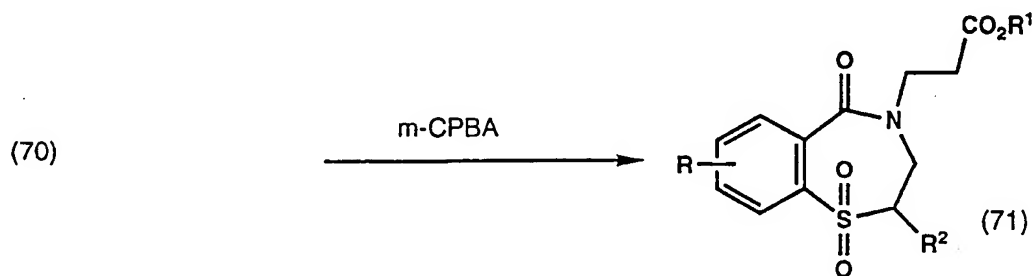




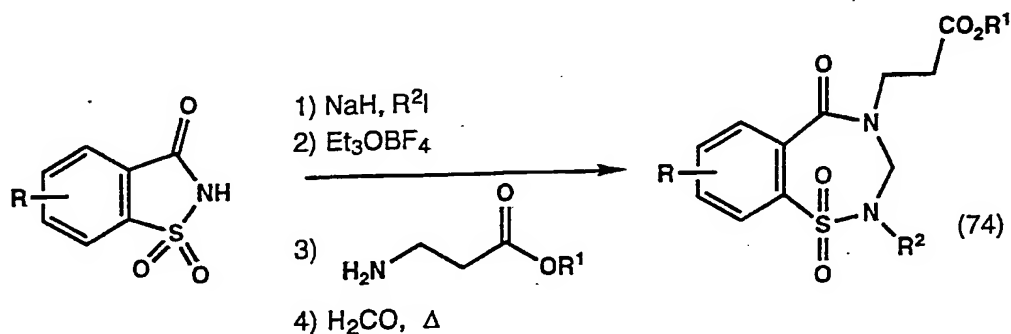
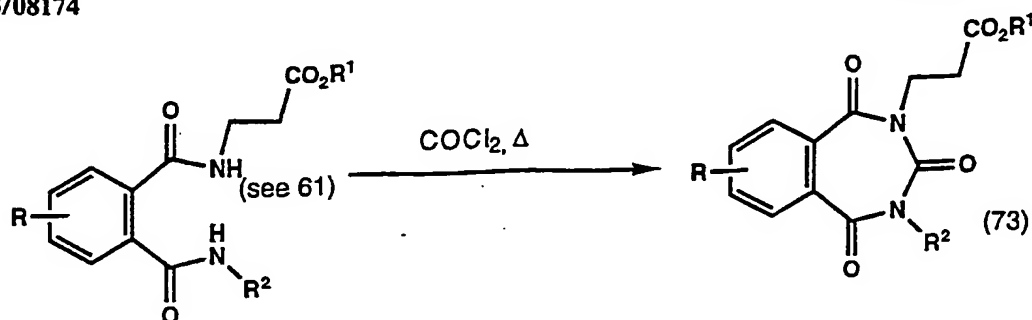
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Starting reagents employed in scheme 8 are either commercially available or readily synthesized by known procedures.

#### 4. Isomeric Products

In products of Formula I carbon atoms bonded to four nonidentical substituents are asymmetric. Accordingly, the compounds may exist as diastereoisomers, enantiomers or mixtures thereof. The syntheses described above may employ racemates, enantiomers or diastereomers as starting materials or intermediates. Diastereomeric products resulting from such syntheses may be separated by chromatographic or crystallization methods. Likewise, enantiomeric product mixtures may be separated using the same techniques or by other methods known in the art. Each of the asymmetric carbon atoms, when present in compounds of Formula I, may be in one of two configurations (R or S) and both are within the scope of the present invention.

#### E. Pharmaceutical Compositions

The compounds described in this invention may be isolated as the free acid or base or converted to salts of various inorganic and organic acids and bases. Such salts are within the scope of this invention. Examples of such salts include ammonium, metal salts like sodium, potassium, calcium and magnesium; salts with organic bases like dicyclohexylamine, N-methyl-D-glucamine and the like; and salts with amino acids like arginine or lysine. Salts with

inorganic and organic acids may be likewise prepared, for example, using hydrochloric, hydrobromic, sulfuric, phosphoric, trifluoroacetic, methanesulfonic, malic, maleic, fumaric and the like. Non-toxic and physiologically compatible salts are particularly useful although other less desirable salts may have use in the processes of isolation and purification.

5           A number of methods are useful for the preparation of the salts described above and are known to those skilled in the art. For example, reaction of the free acid or free base form of a compound of Formula I with one or more molar equivalents of the desired acid or base in a solvent or solvent mixture in which the salt is insoluble; or in a solvent like water after which the solvent is removed by evaporation, distillation or freeze drying. Alternatively, the free acid  
10           or base form of the product may be passed over an ion exchange resin to form the desired salt or one salt form of the product may be converted to another using the same general process.

          In the management of thromboembolic disorders the compounds of this invention may be utilized in compositions such as tablets, capsules or elixirs for oral administration; suppositories for rectal administration; sterile solutions or suspensions for injectable  
15           administration, and the like. Animals in need of treatment using compounds of this invention can be administered dosages that will provide optimal efficacy. The dose and method of administration will vary from animal to animal and be dependent upon such factors as weight, diet, concurrent medication and other factors which those skilled in the medical arts will recognize.

20           Dosage formulations of the nonpeptidyl inhibitors of the present invention are prepared for storage or administration by mixing the inhibitor having the desired degree of purity with physiologically acceptable carriers, excipients, or stabilizers. Such materials are non-toxic to the recipients at the dosages and concentrations employed, and include buffers such as phosphate, citrate, acetate and other organic acid salts; antioxidants such as ascorbic  
25           acid; low molecular weight (less than about ten residues) peptides such as polyarginine, proteins, such as serum albumin, gelatin, or immunoglobulins; hydrophilic polymers such as polyvinylpyrrolidinone; amino acids such as glycine, glutamic acid, aspartic acid, or arginine; monosaccharides, disaccharides, and other carbohydrates including cellulose or its derivatives, glucose, mannose, or dextrans; chelating agents such as EDTA; sugar alcohols such as mannitol  
30           or sorbitol; counterions such as sodium and/or nonionic surfactants such as Tween, Pluronics or polyethyleneglycol.

          Dosage formulations of the nonpeptidyl inhibitors of the present invention to be used for therapeutic administration must be sterile. Sterility is readily accomplished by filtration through sterile filtration membranes such as 0.2 micron membranes. Nonpeptidyl inhibitor  
35           formulations ordinarily will be stored in lyophilized form or as an aqueous solution. The pH of the cyclic inhibitor preparations typically will be between 3 and 11, more preferably from 5 to 9 and most preferably from 7 to 8. It will be understood that use of certain of the foregoing excipients, carriers, or stabilizers will result in the formation of cyclic polypeptide salts. While the preferred route of administration is by hypodermic injection needle, other methods of

administration are also anticipated such as suppositories, aerosols, oral dosage formulations and topical formulations such as ointments, drops and dermal patches.

Therapeutic nonpeptidyl inhibitor formulations generally are placed into a container having a sterile access port, for example, an intravenous solution bag or vial having a stopper pierceable by hypodermic injection needle.

Therapeutically effective dosages may be determined by either *in vitro* or *in vivo* methods. For each particular nonpeptidyl inhibitor of the present invention, individual determinations may be made to determine the optimal dosage required. The range of therapeutically effective dosages will naturally be influenced by the route of administration. For injection by hypodermic needle it may be assumed the dosage is delivered into the body's fluids. For other routes of administration, the absorption efficiency must be individually determined for each inhibitor by methods well known in pharmacology.

The range of therapeutic dosages is from about 0.001 nM to 1.0 mM, more preferably from 0.1 nM to 100 mM, and most preferably from 1.0 nM to 50 mM.

Typical formulation of compounds of Formula I as pharmaceutical compositions are discussed below.

About 0.5 to 500 mg of a compound or mixture of compounds of Formula I, as the free acid or base form or as a pharmaceutically acceptable salt, is compounded with a physiologically acceptable vehicle, carrier, excipient, binder, preservative, stabilizer, flavor, etc., as called for by accepted pharmaceutical practice. The amount of active ingredient in these compositions is such that a suitable dosage in the range indicated is obtained.

Typical adjuvants which may be incorporated into tablets, capsules and the like are a binder such as acacia, corn starch or gelatin; an excipient such as microcrystalline cellulose; a disintegrating agent like corn starch or alginic acid; a lubricant such as magnesium stearate; a sweetening agent such as sucrose or lactose; a flavoring agent such as peppermint, wintergreen or cherry. When the dosage form is a capsule, in addition to the above materials it may also contain a liquid carrier such as a fatty oil. Other materials of various types may be used as coatings or as modifiers of the physical form of the dosage unit. A syrup or elixir may contain the active compound, a sweetener such as sucrose, preservatives like propyl paraben, a coloring agent and a flavoring agent such as cherry. Sterile compositions for injection can be formulated according to conventional pharmaceutical practice. For example, dissolution or suspension of the active compound in a vehicle such as water or naturally occurring vegetable oil like sesame, peanut, or cottonseed oil or a synthetic fatty vehicle like ethyl oleate or the like may be desired. Buffers, preservatives, antioxidants and the like can be incorporated according to accepted pharmaceutical practice.

#### F. Platelet Inhibition Assays

Evaluation of the Formula I inhibitors of the fibrinogen-platelet interaction is guided by *in vitro* receptor binding assays and *in vitro* platelet aggregation inhibition assays.

In-vitro biological activity of the compounds of Formula I are monitored using a modified fibrinogen-GPIIb/IIIa ELISA based on the method of Nachman and Leung (*J. Clin. Invest.* 69:263-269 (1982)) which measures the inhibition of fibrinogen binding to purified human platelet GPIIb/IIIa receptor. Human fibrinogen is prepared by the method of Lipinska, *et al.* (*J. Lab. Clin. Med.* 84:509-516 (1974)). Platelet GPIIb/IIIa is prepared by the method of Fitzgerald, *et al.*, *Ann. Biochem.*, 151:169-177 (1985).

Briefly, microtiter plates are coated with fibrinogen (10 mg/ml) and then blocked with TACTS buffer containing 0.5% bovine serum albumin (BSA). (TACTS buffer contains 20mM Tris.HCl, pH 7.5, 0.02% sodium azide, 2 mM calcium chloride, 0.05% Tween 20, 150 mM sodium chloride.) The plate is washed with phosphate buffered saline (PBS) containing 0.01% Tween 20 and the sample to be determined added, followed by addition of solubilized GPIIb/IIIa receptor (40 mg/ml) in TACTS, 0.5% BSA. After incubation, the plate is washed and 1 mg/ml of murine anti-platelet monoclonal antibody AP3 (Newman *et al.*, *Blood* 65:227-232 (1985)) is added. After another wash a goat anti-mouse IgG conjugated to horseradish peroxidase is added. A final wash is performed and developing reagent buffer (10 mg o-phenylenediamine dihydrochloride, 0.0212% hydrogen peroxide, 0.22 mM citrate, 50 mM phosphate, pH 5.0) is added and then incubated until color develops. The reaction is stopped with 1N sulfuric acid and the absorbance at 492 nm is recorded.

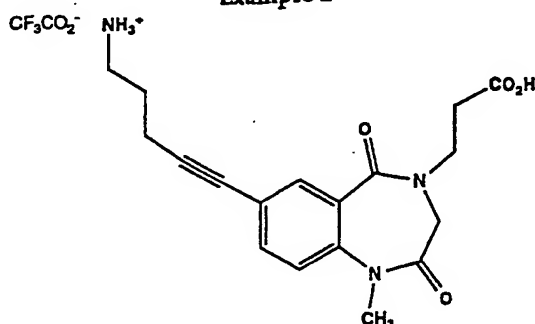
In addition to the GPIIb/IIIa ELISA assay, platelet aggregation assays may be performed in human platelet rich plasma (PRP). Fifty milliliters of whole human blood (9 parts) is drawn on 3.6% sodium citrate (1 part) from a donor who has not taken aspirin or related medications for at least two weeks. The blood is centrifuged at 160 x g for 10 min at 22°C and then allowed to stand for 5 min after which the PRP is decanted. Platelet poor plasma (PPP) is isolated from the remaining blood after centrifugation at 2000 x g for 25 min. The platelet count of the PRP is adjusted to ca. 300,000 per microliter with PPP.

A 225 mL aliquot of PRP plus 25 mL of either a dilution of the test sample or a control (PBS) is incubated for 5 min in a Chrono-log Whole Blood Aggregometer at 25°C. An aggregating agent (collagen, 1 mg/ml; U46619, 100 ng/ml; or ADP, 8 mM) is added and the platelet aggregation recorded.

Without further description, it is believed that one of ordinary skill in the art can, using the preceding description and illustrative examples, make and utilize the present example to the fullest extent. The following working examples therefore, specifically point out preferred embodiments of the present invention, and are not to be construed as limiting in any way the remainder of the disclosure.

## EXAMPLES

## Example 1



5 **1-methyl-4-(2-carboxyethyl)-7-(5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate.**

a) To a mechanically stirred solution of 26.3 grams of 5-iodo-2-amino benzoic acid (0.1 mol), 10.6 grams of sodium carbonate (0.1 mol), and 250 mL water, cooled to 0°C, was slowly added, *via* an addition funnel, 80mL of a 1.93M solution of phosgene in toluene. After 2 hours,  
10 the precipitated product was isolated by filtration. The solids were washed with 200mL water, 300mL of a 1:1 mixture ethanol and ether, 200mL of ether, and dried under vacuum to yield 24.3 grams (84%) of 5-iodoisatoic anhydride (264-268°C, decomposition).

b) A magnetically stirred solution of 5 grams of 5-iodoisatoic anhydride (0.0173 mol),  
15 5.85 grams of b-alanine benzyl ester tosylate (0.0173 mol), 35mL pyridine, and 0.5 of dimethylaminopyridine (0.0041 mol) was heated to 80°C for 2 hrs. The reaction mixture was allowed to cool to room temperature and concentrated *in vacuo*. The resulting residue was dissolved in 100mL ethyl acetate and washed 2X50mL of 10% cupric sulfate, 1X50mL sat. sodium bicarbonate, 1X 50mL brine, dried over sodium sulfate, filtered and concentrated *in*  
20 *vacuo*. The product was further purified by column chromatography, using silica gel, eluting with a 1:1 mixture of ethyl acetate and hexane (TLC, SiO<sub>2</sub>, 1:1 EtOAc/hexane, R<sub>f</sub>=0.65, un positive) to yield 1.85 grams (25%) of N-(2-amino-5-iodobenzoyl)-b-alanine benzyl ester. <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 7.54 (1H, d, <sup>4</sup>J<sub>HH</sub> = 2Hz, Ar-H *o*-CON), 7.42 (1H, dd, <sup>3</sup>J<sub>HH</sub>=9Hz, <sup>4</sup>J<sub>HH</sub>=2Hz, Ar-H *p*-CON), 7.38-7.32 (5H, s, ArH Ph), 6.65 (1H, bt, <sup>3</sup>J<sub>HH</sub>=6Hz, CONH), 6.46  
25 (1H, d, <sup>3</sup>J<sub>HH</sub>=9Hz, Ar-H *m*-CON), 5.16 (2H, s, OCH<sub>2</sub>), 3.68 (2H, q, <sup>3</sup>J<sub>HH</sub>=6Hz, NCH<sub>2</sub>), 2.69 (2H, t, <sup>3</sup>J<sub>HH</sub>=6Hz, CH<sub>2</sub>CO<sub>2</sub>). <sup>13</sup>C NMR (CDCl<sub>3</sub>,dTMS) 172.4, 167.8, 148.2, 140.6, 135.6, 135.5, 128.7, 128.4, 128.3, 119.3, 118.1, 76.3, 66.7, 35.2, 34.0.

Using the above procedure, but substituting the appropriate 3-aminopropionate alkyl ester for b-alanine ethyl ester and N-substituted-5-iodo-isatoic anhydride for 5-  
30 iodoisatoic anhydride there may be prepared, for example, the following compounds:  
ethyl N-(2-amino-5-iodobenzoyl)-3-amino-3-methylpropanoate,  
ethyl N-(2-amino-5-iodobenzoyl)-3-amino-3-phenylpropanoate,

- ethyl N-(2-amino-5-iodobenzoyl)-3-amino-3-(3-(methoxycarbonyl)phenyl)propanoate,  
t-butyl N-[N-(2-amino-5-iodobenzoyl)-L-(b-benzyl)-aspartinate,  
ethyl N-[N-(2-amino-5-iodobenzoyl)-L-aspartyl-(b-benzyl ester)]-glycinate,  
ethyl N-[N-(2-amino-5-iodobenzoyl)-L-aspartyl-(b-benzyl ester)]-valinate,  
5 ethyl N-[N-(2-amino-5-iodobenzoyl)-L-aspartyl-(b-benzyl ester)]-phenylalaninate,  
ethyl N-(2-(N-methyl)-amino-5-iodobenzoyl)-3-amino-3-methylpropanoate,  
ethyl N-(2-(N-methyl)-amino-5-iodobenzoyl)-3-amino-3-phenylpropanoate,  
ethyl N-(2-(N-methyl)-amino-5-iodobenzoyl)-3-amino-3-(3-(methoxycarbonyl)phenyl)propanoate,  
10 t-butyl N-[N-[2-(N-methyl)-amino-5-iodobenzoyl]-(±)-(b-benzyl)-aspartinate,  
ethyl N-[N-(2-(N-methyl)-amino-5-iodobenzoyl)-(±)-aspartyl-(b-benzyl ester)]-glycinate,  
ethyl N-[N-(2-(N-methyl)-amino-5-iodobenzoyl)-(±)-aspartyl-(b-benzyl ester)]-valinate,  
ethyl N-[N-(2-(N-methyl)-amino-5-iodobenzoyl)-(±)-aspartyl-(b-benzyl ester)]-phenylalaninate,  
15 ethyl N-(2-(N-phenyl)-amino-5-iodobenzoyl)-3-amino-3-methylpropanoate,  
ethyl N-(2-(N-phenyl)-amino-5-iodobenzoyl)-3-amino-3-phenylpropanoate,  
ethyl N-(2-(N-phenyl)-amino-5-iodobenzoyl)-3-amino-3-(3-(methoxycarbonyl)phenyl)propanoate,  
t-butyl N-[N-[2-(N-phenyl)-amino-5-iodobenzoyl]-(±)-(b-benzyl)-aspartinate,  
20 ethyl N-[N-(2-(N-phenyl)-amino-5-iodobenzoyl)-(±)-aspartyl-(b-benzyl ester)]-glycinate,  
ethyl N-[N-(2-(N-phenyl)-amino-5-iodobenzoyl)-(±)-aspartyl-(b-benzyl ester)]-valinate,  
ethyl N-[N-(2-(N-phenyl)-amino-5-iodobenzoyl)-(±)-aspartyl-(b-benzyl ester)]-phenylalaninate,  
ethyl N-(2-(N-benzyl)-amino-5-iodobenzoyl)-3-amino-3-methylpropanoate,  
25 ethyl N-(2-(N-benzyl)-amino-5-iodobenzoyl)-3-amino-3-phenylpropanoate,  
ethyl N-(2-(N-benzyl)-amino-5-iodobenzoyl)-3-amino-3-(3-(methoxycarbonyl)phenyl)propanoate,  
t-butyl N-[N-[2-(N-benzyl)-amino-5-iodobenzoyl]-(±)-(b-benzyl)-aspartinate,  
ethyl N-[N-(2-(N-benzyl)-amino-5-iodobenzoyl)-(±)-aspartyl-(b-benzyl ester)]-glycinate,  
30 ethyl N-[N-(2-(N-benzyl)-amino-5-iodobenzoyl)-(±)-aspartyl-(b-benzyl ester)]-valinate,  
ethyl N-[N-(2-(N-benzyl)-amino-5-iodobenzoyl)-(±)-aspartyl-(b-benzyl ester)]-phenylalaninate,  
ethyl N-(2-(N-isopropyl)-amino-5-iodobenzoyl)-3-amino-3-methylpropanoate,  
ethyl N-(2-(N-isopropyl)-amino-5-iodobenzoyl)-3-amino-3-phenylpropanoate,  
35 ethyl N-(2-(N-isopropyl)-amino-5-iodobenzoyl)-3-amino-3-(3-(methoxycarbonyl)phenyl)propanoate,  
t-butyl N-[N-[2-(N-isopropyl)-amino-5-iodobenzoyl]-(±)-(b-benzyl)-aspartinate,  
ethyl N-[N-(2-(N-isopropyl)-amino-5-iodobenzoyl)-(±)-aspartyl-(b-benzyl ester)]-glycinate,

ethyl N-[N-(2-(N-isopropyl)-amino-5-iodobenzoyl)-(±)-aspartyl-(b-benzyl ester)]-  
valinate,  
ethyl N-[N-(2-(N-isopropyl)-amino-5-iodobenzoyl)-(±)-aspartyl-(b-benzyl ester)]-  
phenylalaninate,

5

c) A magnetically stirred solution of 0.848 gram of N-(2-amino-5-iodobenzoyl)-b-  
alanine ethyl ester (2.0 mmol), 0.35 mL 2,6-lutidine (3.0 mmol), 0.19 mL methyl iodide (3.0  
mmol), and 15 mL dimethylformamide was heated to 50°C for 15 hours. The reaction mixture  
was allowed to cool to room temperature and concentrated *in vacuo*. The resulting residue was  
10 dissolved in 75 mL ethyl acetate and washed 1X50 mL 10% citric acid, 1X50 mL sat. sodium  
bicarbonate, 1X50 mL brine, dried over magnesium sulfate, filtered and concentrated *in vacuo*.  
The resulting oil was further purified by column chromatography, using silica gel, eluting  
with a solvent gradient of 35/65 ethyl acetate/hexane to 65/35 ethyl acetate/hexane (TLC,  
SiO<sub>2</sub>, 1:1 EtOAc/hexane, R<sub>f</sub>=0.84, un positive) to yield 305 mgs (35%) of N-(2-methylamino-  
15 5-iodobenzoyl)-b-alanine benzyl ester. <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 7.56 (2H, m, Ar-H *o,p*-  
CON), 7.43 (1H, bq, <sup>3</sup>J<sub>HH</sub>=5Hz, NHMe), 7.38-7.32 (5H, s, ArH Ph), 6.62 (1H, bt, <sup>3</sup>J<sub>HH</sub>=6Hz,  
CONH), 6.42 (1H, d, <sup>3</sup>J<sub>HH</sub>=9Hz, Ar-H *m*-CON), 5.16 (2H, s, OCH<sub>2</sub>), 3.64 (2H, q, <sup>3</sup>J<sub>HH</sub>=6Hz,  
NCH<sub>2</sub>), 2.81 (3H, d, <sup>3</sup>J<sub>HH</sub>=5Hz, NCH<sub>3</sub>), 2.64 (2H, t, <sup>3</sup>J<sub>HH</sub>=6Hz, CH<sub>2</sub>CO<sub>2</sub>). <sup>13</sup>C NMR  
(CDCl<sub>3</sub>, dTMS) 172.4, 168.4, 148.8, 141.0, 135.5, 128.7, 128.4, 128.3, 117.3, 113.4, 74.1, 66.7, 35.2,  
20 34.0, 29.6.

d) To a magnetically stirred solution of 0.305 grms of N-(2-methylamino-5-  
iodobenzoyl)-b-alanine benzyl ester (0.69 mmol), 3 mL methylene chloride, and 0.144 mL  
triethylamine (1.04 mmol), cooled to -30°C under an atmosphere of nitrogen was slowly added  
25 0.09 mL of α-bromoacetyl bromide (1.04 mmol) as a solution in 2 mL methylene chloride. The  
reaction mixture was allowed to warm to room temperature and stir for 2 hours. The mixture  
was diluted with 40 mL methylene chloride and washed with 1X50 10% citric acid, 1X50 sat.  
sodium bicarbonate, dried over sodium sulfate, filtered and concentrated *in vacuo*. The  
resulting residue was dissolved in 3 mL dimethylformamide and added, *via* an addition  
30 funnel, to a slurry of 25 mgs sodium hydride (1.04 mmol) in 2 mL dimethylformamide that was  
cooled to 0°C. After 2 hours, the mixture was poured over 50 mL of an ice cooled solution of 10%  
citric acid and extracted with 3X40 mL ethyl acetate. The combined organic layers were  
washed with 1X50 mL sat. sodium bicarbonate, dried over magnesium sulfate, and  
concentrated *in vacuo*. The resulting residue was further purified by column chromatography,  
35 using silica gel, eluting with a solvent gradient of 40:60 ethyl acetate/hexane to 70:30 ethyl  
acetate/hexane to yield 0.16 gms (49%) of 1-methyl-4-(2-carboxyethyl)-7-iodo-3,4-dihydro-  
1H-1,4-benzodiazepine-2,5-dione benzyl ester (TLC, SiO<sub>2</sub>, 1:1 ethyl acetate/hexane,  
R<sub>f</sub>=0.48, un positive). <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 8.15 (1H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, Ar-H *o*-CON), 7.78  
(1H, dd, <sup>4</sup>J<sub>HH</sub>=2Hz, <sup>3</sup>J<sub>HH</sub>=9Hz, Ar-H *p*-CON), 6.90 (1H, d, <sup>3</sup>J<sub>HH</sub>=9Hz, Ar-H *m*-CON), 5.1



(2H, s, OCH<sub>2</sub>), 3.92 (1H, d, <sup>2</sup>J<sub>HH</sub>=14Hz, COCHHN), 3.90 (2H, t, <sup>3</sup>J<sub>HH</sub>=8Hz, NCH<sub>2</sub>), 3.82 (1H, d, <sup>2</sup>J<sub>HH</sub>=14Hz, COCHHN) 3.27 (3H, s, NCH<sub>3</sub>), 2.82 (1H, dt, <sup>2</sup>J<sub>HH</sub>=19Hz, <sup>3</sup>J<sub>HH</sub>=9Hz, CHHCO<sub>2</sub>), 2.68 (1H, dt, <sup>2</sup>J<sub>HH</sub>=19Hz, <sup>3</sup>J<sub>HH</sub>=9Hz, CHHCO<sub>2</sub>).

Using the above procedure, but substituting the appropriate N-(2-amino-5-

- 5 iodobenzoyl)-b-alanine alkyl ester for N-(2-methylamino-5-iodobenzoyl)-b-alanine benzyl ester and a-substituted-a-halo acetyl halide for a-bromo acetyl bromide there may be prepared, for example, the following compounds:  
1-methyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 10 1-methyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl)]-7-iodo-3,4-dihydro-1H-1,4-
- 15 benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-benzyloxyvaliny-4-benzyloxy)succinyl)]-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-benzyloxyphenylalany-4-benzyloxy)succinyl)]-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 20 1-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-iodo-3,4-dihydro-1H-1,4-
- 25 benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl)]-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyvaliny-4-benzyloxy)succinyl)]-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 30 (±)-1-phenyl-4-[2-(1-benzyloxyphenylalany-4-benzyloxy)succinyl)]-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-benzyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-benzyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-iodo-3,4-dihydro-1H-1,4-
- 35 benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl)]-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-benzyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 1-isopropyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-isopropyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 (±)-1-isopropyl-4-[2-(1-benzyloxyglycinyl-4-benzyloxy)succinyl)]-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 (±)-1-isopropyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-methyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-methyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 (±)-1-methyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyglycinyl-4-benzyloxy)succinyl)]-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 (±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-iodo-3,4-dihydro-3-phenyl-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1,3-diphenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 1,3-diphenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 (±)-1,3-diphenyl-4-[2-(1-benzyloxyglycinyl-4-benzyloxy)succinyl)]-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1,3-diphenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-benzyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 5 1-benzyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl)]-7-iodo-3,4-
- 10 dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 15 1-isopropyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-isopropyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-iodo-3,4-dihydro-1H-
- 20 1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl)]-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 25 (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

e) To a magnetically stirred solution of 9.13 grams hex-5-ynoic acid (0.081 mol), 7.12mL triethylamine (0.051 mol), 30mL *tert*-butanol, was added 17.7mL diphenylphosphoryl azide (0.082 mol) and the mixture heated to reflux for 3 hours. The reaction mixture was allowed to cool to room temperature and poured over a biphasic mixture of 200mL water and 200mL ether. The layers were separated and the organic layer was washed 1X50mL 5% ethylene-

30 diaminetetraacetic acid disodium salt, 1X50mL 10% sodium bicarbonate, 1X50mL brine, dried over magnesium sulfate and activated carbon, filtered and concentrated *in vacuo*. The

35 resulting residue was further purified by column chromatography, using silica gel, eluting with 1:1 ether/hexane (TLC, SiO<sub>2</sub>, 1:3 ethyl acetate/hexane, R<sub>f</sub>=0.18, ninhydrin char) to yield 3.8 grams (25%) N-boc-5-amino-1-pentyne. <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 4.65 (1H, bs, NH), 3.22 (2H, q, <sup>3</sup>J<sub>HH</sub>=6Hz, NCH<sub>2</sub>), 2.23 (2H, dt, <sup>3</sup>J<sub>HH</sub>=6Hz, <sup>4</sup>J<sub>HH</sub>=3Hz, CH<sub>2</sub>C(C)), 1.98 (1H, t, <sup>4</sup>J<sub>HH</sub>=3Hz, C(CH<sub>3</sub>)), 1.73 (2H, p, <sup>3</sup>J<sub>HH</sub>=6Hz, CH<sub>2</sub>), 1.42 (9H, s, *t*-Bu)

f) To a magnetically stirred solution of 160mgs 1-methyl-2-(carboxyethyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione benzyl ester (0.335mmol) in 5mL ethyl acetate, degassed of oxygen, under an atmosphere of nitrogen was added 123mgs N-boc-5-amino-1-pentyne (0.67mmol), 10mgs bis-triphenylphosphine palladium dichloride (0.014mmol), 5mgs cuprous iodide (0.026mmol) and 0.233mL triethylamine (1.675mmol). After 2 hours, the reaction mixture was diluted with 50mL ethyl acetate and washed 2X50mL 10% citric acid, 2X50mL sat. sodium bicarbonate, dried over magnesium sulfate, filtered, and concentrated *in vacuo*. The resulting residue was further purified by column chromatography, using silica gel, eluting with a solvent gradient of 50:50 ethyl acetate/hexane to 75:25 ethyl acetate/hexane (TLC, SiO<sub>2</sub>, 1:1 ethyl acetate/hexane, R<sub>f</sub>=0.39, un positive) to yield 90mgs (50%) 1-methyl-4-(2-carboxyethyl)-7-(N-boc-5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester. <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 7.85 (1H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, Ar-H *o*-CON), 7.48 (1H, dd, <sup>4</sup>J<sub>HH</sub>=2Hz, <sup>3</sup>J<sub>HH</sub>=8Hz, Ar-H *p*-CON), 7.32 (5H, s, Ph), 7.06 (1H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, Ar-H *m*-CON), 5.10 (2H, s, OCH<sub>2</sub>), 4.67 (1H, bs, NH), 3.95 (1H, d, <sup>2</sup>J<sub>HH</sub>=15Hz, NCHHCO), 3.91 (2H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, NCH<sub>2</sub>), 3.81 (1H, d, <sup>2</sup>J<sub>HH</sub>=15Hz, NCHHCO), 3.29 (3H, s, NCH<sub>3</sub>), 3.26 (2H, q, <sup>3</sup>J<sub>HH</sub>=7Hz, BocNHCH<sub>2</sub>), 2.82 (1H, dt, <sup>2</sup>J<sub>HH</sub>=16Hz, <sup>3</sup>J<sub>HH</sub>=8Hz, CHHCO<sub>2</sub>), 2.69 (1H, dt, <sup>2</sup>J<sub>HH</sub>=16Hz, <sup>3</sup>J<sub>HH</sub>=8Hz, CHHCO<sub>2</sub>), 2.46 (2H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, C[CCH<sub>2</sub>), 1.78 (2H, p, <sup>3</sup>J<sub>HH</sub>=7Hz, CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>).

Using the above procedure, but substituting the appropriate 7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione for 1-methyl-2-(carboxyethyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione benzyl ester and alkyne for N-boc-5-amino-1-pentyne there may be prepared, for example, the following compounds:

1-methyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 1-methyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 (±)-1-methyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 (±)-1-methyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 (±)-1-methyl-4-[2-(1-benzyloxyvaliny-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 (±)-1-methyl-4-[2-(1-benzyloxyphenylalany-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 1-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 1-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 5 (±)-1-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 1-benzyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 10 1-benzyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-benzyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 15 (±)-1-benzyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-benzyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-benzyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 20 1-isopropyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 1-isopropyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 25 (±)-1-isopropyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-isopropyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-isopropyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 30 (±)-1-isopropyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 1-methyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 35 1-methyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-methyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-methyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl)]-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 5 (±)-1-methyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 10 (±)-1-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl)]-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 15 (±)-1-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-benzyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 20 1-benzyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 25 (±)-1-benzyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl)]-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 30 1-isopropyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-isopropyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 35 (±)-1-isopropyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl)]-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-isopropyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 5 1-methyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-methyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 10 (±)-1-methyl-4-[2-(1-benzyloxyglycinyl-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 15 (±)-1-methyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 20 (±)-1-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyglycinyl-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 25 (±)-1-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-benzyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 30 1-benzyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 35 (±)-1-benzyl-4-[2-(1-benzyloxyglycinyl-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-benzyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-isopropyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 1-isopropyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl)]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 (±)-1-isopropyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 1-methyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-methyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 (±)-1-methyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl)]-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 (±)-1-methyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 (±)-1-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl)]-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 (±)-1-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,



- 1-benzyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-benzyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 5 (±)-1-benzyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl])-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl])-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl])-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 10 (±)-1-benzyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl])-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-isopropyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 15 1-isopropyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl])-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl])-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 20 (±)-1-isopropyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl])-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl])-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 25 1-methyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-methyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl])-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 30 (±)-1-methyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl])-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl])-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 35 (±)-1-methyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl])-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- 1-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 (±)-1-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 (±)-1-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-benzyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-benzyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 (±)-1-benzyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 (±)-1-benzyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-isopropyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-isopropyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 (±)-1-isopropyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 (±)-1-isopropyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 1-methyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperizine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-methyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperizine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-methyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 (±)-1-methyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 1-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 (±)-1-phenyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 1-benzyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-benzyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 (±)-1-benzyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 (±)-1-benzyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-isopropyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 1-isopropyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-isopropyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 (±)-1-isopropyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-methyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-methyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 (±)-1-methyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 (±)-1-methyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 1-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 (±)-1-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 1-benzyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-benzyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 (±)-1-benzyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-benzyl-4-[2-(1-benzyl-oxyvalinyl-4-benzyl-oxy)succinyl)]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 (±)-1-benzyl-4-[2-(1-benzyl-oxyphenylalanyl-4-benzyl-oxy)succinyl)]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 5 1-isopropyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 1-isopropyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 (±)-1-isopropyl-4-[2-(1-*t*-butoxy-4-benzyl-oxy)succinyl)]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 10 (±)-1-isopropyl-4-[2-(1-benzyl-oxyglycyl-4-benzyl-oxy)succinyl)]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 (±)-1-isopropyl-4-[2-(1-benzyl-oxyvalinyl-4-benzyl-oxy)succinyl)]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 15 (±)-1-isopropyl-4-[2-(1-benzyl-oxyphenylalanyl-4-benzyl-oxy)succinyl)]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 1-methyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 1-methyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 20 (±)-1-methyl-4-[2-(1-*t*-butoxy-4-benzyl-oxy)succinyl)]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 (±)-1-methyl-4-[2-(1-benzyl-oxyglycyl-4-benzyl-oxy)succinyl)]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 25 (±)-1-methyl-4-[2-(1-benzyl-oxyvalinyl-4-benzyl-oxy)succinyl)]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 (±)-1-methyl-4-[2-(1-benzyl-oxyphenylalanyl-4-benzyl-oxy)succinyl)]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 1-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 30 (±)-1-phenyl-4-[2-(1-*t*-butoxy-4-benzyl-oxy)succinyl)]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 (±)-1-phenyl-4-[2-(1-benzyl-oxyglycyl-4-benzyl-oxy)succinyl)]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 35 (±)-1-phenyl-4-[2-(1-benzyl-oxyvalinyl-4-benzyl-oxy)succinyl)]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-benzyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 1-benzyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl)]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 (±)-1-benzyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 1-isopropyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-isopropyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 (±)-1-isopropyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl)]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 (±)-1-isopropyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-methyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-methyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 (±)-1-methyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 (±)-1-methyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- 1-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 (±)-1-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 (±)-1-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-benzyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 1-benzyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 (±)-1-benzyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 1-isopropyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-isopropyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 (±)-1-isopropyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 (±)-1-isopropyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-methyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- 1-methyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl])-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 (±)-1-methyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl])-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl])-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl])-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 1-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 (±)-1-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl])-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl])-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl])-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 (±)-1-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl])-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-benzyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 1-benzyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl])-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl])-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 (±)-1-benzyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl])-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl])-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 1-isopropyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-isopropyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,



- (±)-1-isopropyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 (±)-1-isopropyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 (±)-1-methyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 (±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 (±)-1,3-diphenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 (±)-1,3-diphenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 (±)-1,3-diphenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 (±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 (±)-1-isopropyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 (±)-1-methyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 (±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 (±)-1,3-diphenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 (±)-1,3-diphenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1,3-diphenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 5 (±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[5-cyano-1-pentyne]-3,4-
- 10 dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyglycinyl-4-benzyloxy)succinyl)]-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 15 (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyne]-3,4-
- 20 dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyglycinyl-4-benzyloxy)succinyl)]-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 25 (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[5-cyano-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-
- 30 ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 35 (±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyglycinyl-4-benzyloxy)succinyl)]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 (±)-1,3-diphenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 (±)-1,3-diphenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 (±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 (±)-1-isopropyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]]-7-[2-(3-cyanophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-methyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-methyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 5 (±)-1-methyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl])-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyglycinyl-4-benzyloxy)succinyl])-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl])-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 10 (±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl])-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1,3-diphenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1,3-diphenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 15 (±)-1,3-diphenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl])-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1,3-diphenyl-4-[2-(1-benzyloxyglycinyl-4-benzyloxy)succinyl])-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 20 (±)-1,3-diphenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl])-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1,3-diphenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl])-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 25 (±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-benzyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl])-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 30 (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyglycinyl-4-benzyloxy)succinyl])-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl])-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 35 (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl])-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[2-(4-nitrophenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 (±)-1-methyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 (±)-1-methyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 (±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 (±)-1,3-diphenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 (±)-1,3-diphenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 (±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-benzyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 5 (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 10 (±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-isopropyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 15 (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 20 (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-methyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-methyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-methyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 30 (±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 35 (±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1,3-diphenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1,3-diphenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperizine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperizine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 (±)-1,3-diphenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperizine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperizine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperizine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 (±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperizine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperizine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 (±)-1-benzyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperizine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperizine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperizine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperizine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperizine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 (±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperizine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperizine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperizine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperizine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyn-3-[4-(N-Boc)-1-piperizine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 (±)-1-methyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,



- (±)-1-methyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl)]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 (±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 (±)-1,3-diphenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 (±)-1,3-diphenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl)]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 (±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 (±)-1-benzyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl)]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 (±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[2-(N-Boc)aminoethanethiol]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 (±)-1-methyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 (±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 (±)-1,3-diphenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 (±)-1,3-diphenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 (±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 (±)-1-benzyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 (±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyglycinyl-4-benzyloxy)succinyl)]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[2-(4-cyanothiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 (±)-1-methyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyglycinyl-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 (±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 (±)-1,3-diphenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 (±)-1,3-diphenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-benzyloxyglycinyl-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1,3-diphenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 (±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 (±)-1-benzyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyglycinyl-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 (±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 (±)-1-isopropyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyglycinyl-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 (±)-1-methyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-methyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 (±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 (±)-1,3-diphenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 (±)-1,3-diphenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 (±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 (±)-1-benzyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 (±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 5 (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- g) To a solution of 22mgs 1-methyl-4-(2-carboxyethyl)-7-(N-boc-5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione benzyl ester (0.0412mmol) in 0.5mL ethyl acetate was added 3mL sat. HCl in ethyl acetate. After 1 hour, the mixture was concentrated *in vacuo* and diluted with 2mL methanol. To the methanolic solution was added 0.5mL 2N sodium hydroxide. After 1 hour, the reaction was quenched with 3mL acetic acid, concentrated *in vacuo*, diluted with 3mL water and purified by high pressure liquid chromatography, using a 1/2" C-18 reverse-phase column, eluting with a solvent gradient of 10:90 methanol(0.1% trifluoroacetic acid)/ water (0.1% trifluoroacetic acid), time 0 to 10 min, to 50:50 methanol(0.1% trifluoroacetic acid)/water (0.1% trifluoroacetic acid), time 10 min to 40 min, flow = 10ml/min ( $R_t=33.1$  min, un detection 254nm) to yield 13mgs (92%) 1-methyl-4-(2-carboxyethyl)-7-(5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate. HRMS (FAB) molecular ion  $m/z=344.1626$  (calcd.  $C_{18}H_{22}N_3O_4$ , 344.1610).  
 $^1H$  NMR ( $D_2O$ ) 7.63 (1H, d,  $^4J_{HH}$ , Ar-H *o*-CON), 7.54 (1H, dd,  $^4J_{HH}=2Hz$ ,  $^3J_{HH}=8Hz$ , Ar-H *p*-CON), 7.20 (1H, d,  $^3J_{HH}=8Hz$ , Ar-H *m*-CON), 4.05 (1H, d,  $^2J_{HH}=14Hz$ , NCHHCO), 4.0 (1H, m, NCHHCH<sub>2</sub>), 3.74 (1H, d,  $^2J_{HH}=14Hz$ , NCHHCO), 3.59 (1H, m, NCHHCH<sub>2</sub>), 3.20 (3H, s, NCH<sub>3</sub>), 3.03 (2H, t,  $^3J_{HH}=7Hz$ , NCH<sub>2</sub>), 2.59 (2H, m, CH<sub>2</sub>CO<sub>2</sub>), 2.42 (2H, t,  $^3J_{HH}=7Hz$ , C[CCH<sub>2</sub>]), 1.82 (2H, p,  $^3J_{HH}=7Hz$ , CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>).  
 25

Using the above procedure, but substituting the appropriate 3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione for 1-methyl-4-(2-carboxyethyl)-7-(N-boc-5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione benzyl ester there may be prepared, for example, the following compounds:

- 30 1-methyl-4-(2-carboxy-1-methylethyl)-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 1-methyl-4-(2-carboxy-1-phenylethyl)-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 (±)-1-methyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 35 (±)-1-methyl-4-[succin-2-yl-(1-glycine)]-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 (±)-1-methyl-4-[succin-2-yl-(1-valine)]-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1-methyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-phenyl-4-(2-carboxy-1-methylethyl)-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 1-phenyl-4-(2-carboxy-1-phenylethyl)-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-glycine)]-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 (±)-1-phenyl-4-[succin-2-yl-(1-valine)]-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 1-benzyl-4-(2-carboxy-1-methylethyl)-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-benzyl-4-(2-carboxy-1-phenylethyl)-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 (±)-1-benzyl-4-[succin-2-yl-(1-glycine)]-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-valine)]-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1-benzyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-isopropyl-4-(2-carboxy-1-methylethyl)-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-isopropyl-4-(2-carboxy-1-phenylethyl)-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 (±)-1-isopropyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-glycine)]-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-isopropyl-4-[succin-2-yl-(1-valine)]-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- 1-methyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-  
3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-methyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-  
3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
5 (±)-1-methyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-  
3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-  
3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-  
10 3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-  
amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-phenyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-  
3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
15 1-phenyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-  
3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-  
3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-  
20 3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-  
3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-  
amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
25 1-benzyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-  
3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-benzyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-  
3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-  
30 3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-  
3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-  
3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
35 (±)-1-benzyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-  
amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-isopropyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-  
3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,



- 1-isopropyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1-isopropyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-
- 10 3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-methyl-4-(2-carboxy-1-methylethyl)-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-methyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1-methyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-glycine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-valine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-
- 20 1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-phenyl-4-(2-carboxy-1-methylethyl)-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 1-phenyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-glycine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-
- 30 1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-valine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 1-benzyl-4-(2-carboxy-1-methylethyl)-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-benzyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1-benzyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-4-[succin-2-yl-(1-glycine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1-benzyl-4-[succin-2-yl-(1-valine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-isopropyl-4-(2-carboxy-1-methylethyl)-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 1-isopropyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1-isopropyl-4-[succin-2-yl-(1-glycine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-4-[succin-2-yl-(1-valine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 1-methyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-methyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1-methyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 1-phenyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-phenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1-phenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-phenyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-benzyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-benzyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 (±)-1-benzyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1-benzyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-isopropyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 1-isopropyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1-isopropyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 1-methyl-4-(2-carboxy-1-methylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-methyl-4-(2-carboxy-1-phenylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-methyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-4-[succin-2-yl-(1-glycine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1-methyl-4-[succin-2-yl-(1-valine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-phenylalanine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
5 1-phenyl-4-(2-carboxy-1-methylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-phenyl-4-(2-carboxy-1-phenylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
10 (±)-1-phenyl-4-[succin-2-yl-(1-glycine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-valine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
15 (±)-1-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-benzyl-4-(2-carboxy-1-methylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-benzyl-4-(2-carboxy-1-phenylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
20 (±)-1-benzyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-glycine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
25 (±)-1-benzyl-4-[succin-2-yl-(1-valine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-phenylalanine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-isopropyl-4-(2-carboxy-1-methylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
30 1-isopropyl-4-(2-carboxy-1-phenylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-butyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
35 (±)-1-isopropyl-4-[succin-2-yl-(1-glycine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)]-4-oxa-1-butyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-valine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1-isopropyl-4-[succin-2-yl-(1-phenylalanine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-methyl-4-(2-carboxy-1-methylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 1-methyl-4-(2-carboxy-1-phenylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-glycine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 (±)-1-methyl-4-[succin-2-yl-(1-valine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-phenylalanine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 1-phenyl-4-(2-carboxy-1-methylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-phenyl-4-(2-carboxy-1-phenylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 (±)-1-phenyl-4-[succin-2-yl-(1-glycine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-valine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-benzyl-4-(2-carboxy-1-methylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 30 1-benzyl-4-(2-carboxy-1-phenylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-glycine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-benzyl-4-[succin-2-yl-(1-valine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-phenylalanine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- 1-isopropyl-4-(2-carboxy-1-methylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-  
3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-isopropyl-4-(2-carboxy-1-phenylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-  
3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
5 (±)-1-isopropyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-  
3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-glycine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-  
3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-valine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-  
10 3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-phenylalanine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-  
pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-(5-amino-1-pentyne)-3,4-dihydro-  
1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
15 (±)-1-methyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-(5-amino-1-pentyne)-3,4-dihydro-  
1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-amino-1-pentyne)-3,4-dihydro-1H-  
1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-(5-amino-1-pentyne)-3,4-dihydro-1H-  
20 1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-(5-amino-1-pentyne)-3,4-dihydro-1H-  
1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-amino-1-pentyne)-3,4-  
dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
25 (±)-1,3-diphenyl-4-(2-carboxy-1-methylethyl)-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-  
benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-(2-carboxy-1-phenylethyl)-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-  
benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-  
30 benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-[succin-2-yl-(1-glycine)]-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-  
benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-[succin-2-yl-(1-valine)]-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-  
benzodiazepine-2,5-dione trifluoroacetate,  
35 (±)-1,3-diphenyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-amino-1-pentyne)-3,4-dihydro-1H-  
1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-(5-amino-1-pentyne)-3,4-dihydro-  
1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-amino-1-pentyne)-3,4-
- 10 dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-(5-amino-1-pentyne)-3,4-dihydro-
- 20 1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1-methyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-
- 30 3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1,3-diphenyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1,3-diphenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1,3-diphenyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[3-keto-1-propyn-3-(1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-methyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyn-3-(1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyn-3-(1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,



- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-(2-carboxy-1-methylethyl)-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 (±)-1,3-diphenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-glycine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1,3-diphenyl-4-[succin-2-yl-(1-valine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

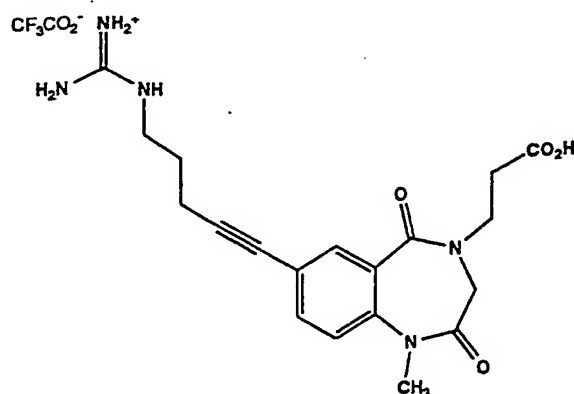
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[3-keto-1-propyn-3-(1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyn-3-(1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1-methyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 (±)-1,3-diphenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1,3-diphenyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(2-aminoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1-methyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-(2-carboxy-1-methylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 (±)-1,3-diphenyl-4-(2-carboxy-1-phenylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-glycine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1,3-diphenyl-4-[succin-2-yl-(1-valine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1-methyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1,3-diphenyl-4-(2-carboxy-1-methylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-(2-carboxy-1-phenylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1,3-diphenyl-4-[succin-2-yl-(1-t-butoxy)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-[succin-2-yl-(1-glycine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-[succin-2-yl-(1-valine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-
- 10 3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 15 (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-t-butoxy)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-
- 20 pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[5-(2-aminopyridine-3-yl)-5-
- 25 oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-t-butoxy)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-
- 30 pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-
- 35 pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate.

## Example 2



5 **1-methyl-4-(2-carboxyethyl)-7-(5-guanidino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate.**

To a magnetically stirred solution of 6mgs of 1-methyl-4-(2-carboxyethyl)-7-(5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate (0.017mmol) in 1.0mL 10% potassium bicarbonate was added 31.0mgs of formamidine sulfonic acid (0.25 mmol). After 30 minutes, the reaction mixture was quenched with 0.5mL acetic acid and concentrated *in vacuo*. The resulting residue was diluted with 2mL water and purified by high pressure liquid chromatography, using a 1/2" C-18 reverse-phase column, eluting with a solvent gradient of 10:90 methanol(0.1% trifluoroacetic acid)/ water (0.1% trifluoroacetic acid), time 0 to 10 min, to 50:50 methanol(0.1% trifluoroacetic acid)/water (0.1% trifluoroacetic acid), time 10 min to 40 min, flow=10ml/min ( $R_t$ =36.8 min, un detection 254nm) to yield 4.5mgs (67%) 1-methyl-4-(2-carboxyethyl)-7-(5-guanidino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate. HRMS (FAB) molecular ion  $m/z$ =386.1823 ( cald.  $C_{19}H_{24}N_5O_4$ , 386.1829).  $^1H$  NMR ( $D_2O$ ) 7.60 (1H, d,  $^4J_{HH}$ , Ar-H *o*-CON), 7.51 (1H, dd,  $^4J_{HH}$ =2Hz,  $^3J_{HH}$ =8Hz, Ar-H *p*-CON), 7.20 (1H, d,  $^3J_{HH}$ =8Hz, Ar-H *m*-CON), 4.05 (1H, d,  $^2J_{HH}$ =14Hz, NCHHCO), 3.92 (1H, m, NCHHCH<sub>2</sub>), 3.72 (1H, d,  $^2J_{HH}$ =14Hz, NCHHCO), 3.40 (1H, m, NCHHCH<sub>2</sub>), 3.23-3.15 (5H, NCH<sub>3</sub>, H<sub>2</sub>NC(=NH<sub>2</sub>)NHCH<sub>2</sub>), 2.58 (2H, m, CH<sub>2</sub>CO<sub>2</sub>), 2.38 (2H, t,  $^3J_{HH}$ =7Hz, C(CCH<sub>2</sub>), 1.76 (2H, p,  $^3J_{HH}$ =7Hz, CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>).

Using the above procedure, but substituting the appropriate amino acid for 1-methyl-4-(2-carboxyethyl)-7-(5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate there may be prepared, for example, the following compounds:

25 1-methyl-4-(2-carboxy-1-methylethyl)-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 1-methyl-4-(2-carboxy-1-phenylethyl)-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 30 (±)-1-methyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1-methyl-4-[succin-2-yl-(1-glycine)]-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-valine)]-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1-methyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-phenyl-4-(2-carboxy-1-methylethyl)-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-phenyl-4-(2-carboxy-1-phenylethyl)-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-
- 10 benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-glycine)]-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1-phenyl-4-[succin-2-yl-(1-valine)]-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-benzyl-4-(2-carboxy-1-methylethyl)-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-
- 20 benzodiazepine-2,5-dione trifluoroacetate,  
1-benzyl-4-(2-carboxy-1-phenylethyl)-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1-benzyl-4-[succin-2-yl-(1-glycine)]-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-valine)]-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-
- 30 1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-isopropyl-4-(2-carboxy-1-methylethyl)-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-isopropyl-4-(2-carboxy-1-phenylethyl)-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-isopropyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-glycine)]-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1-isopropyl-4-[succin-2-yl-(1-valine)]-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 1-methyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-methyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 (±)-1-methyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1-methyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-phenyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-phenyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 (±)-1-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1-phenyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 1-benzyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-benzyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-benzyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,



- (±)-1-benzyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-isopropyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 1-isopropyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 (±)-1-isopropyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,...
- 15 1-methyl-4-(2-carboxy-1-methylethyl)-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-methyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 (±)-1-methyl-4-[succin-2-yl-(1-glycine)]-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-valine)]-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1-methyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-phenyl-4-(2-carboxy-1-methylethyl)-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-phenyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 (±)-1-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-glycine)]-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-phenyl-4-[succin-2-yl-(1-valine)]-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- 1-benzyl-4-(2-carboxy-1-methylethyl)-7-[3-keto-1-propyn-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-benzyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyn-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1-benzyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyn-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-glycine)]-7-[3-keto-1-propyn-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-valine)]-7-[3-keto-1-propyn-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 (±)-1-benzyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyn-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-isopropyl-4-(2-carboxy-1-methylethyl)-7-[3-keto-1-propyn-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 1-isopropyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyn-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyn-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-glycine)]-7-[3-keto-1-propyn-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 (±)-1-isopropyl-4-[succin-2-yl-(1-valine)]-7-[3-keto-1-propyn-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyn-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 1-methyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-methyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 (±)-1-methyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-methyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-phenyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- 1-phenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1-phenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-
- 10 3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-benzyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-benzyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1-benzyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-
- 20 dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-isopropyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 1-isopropyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-
- 30 dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-
- 35 3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
5 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-(2-carboxy-1-methylethyl)-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
10 (±)-1,3-diphenyl-4-(2-carboxy-1-phenylethyl)-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
15 (±)-1,3-diphenyl-4-[succin-2-yl-(1-glycine)]-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-[succin-2-yl-(1-valine)]-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
20 (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
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25 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
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30 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
35 (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 (±)-1,3-diphenyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1,3-diphenyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

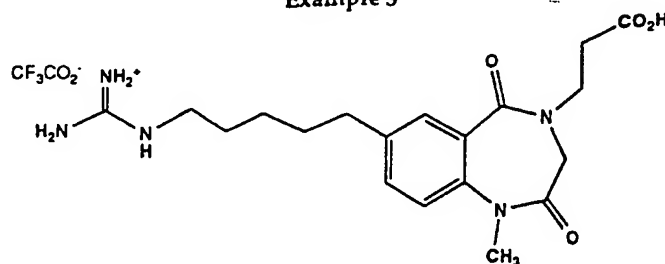
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
5 (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
10 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
15 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
20 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
25 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-(2-carboxy-1-methylethyl)-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
30 (±)-1,3-diphenyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
35 (±)-1,3-diphenyl-4-[succin-2-yl-(1-glycine)]-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-[succin-2-yl-(1-valine)]-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1,3-diphenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 5 (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
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 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
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 (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 20 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 25 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyn-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 (±)-1-methyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 (±)-1-methyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 30 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 35 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1,3-diphenyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
5 (±)-1,3-diphenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
10 (±)-1,3-diphenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
15 (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
20 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
25 (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
30 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
35 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(2-guanidinoethanethiol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate.



Example 3



**1-methyl-4-(2-carboxyethyl)-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate.**

5

- a) A magnetically stirred slurry of 45mgs 1-methyl-4-(2-carboxyethyl)-7-(N-boc-5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione benzyl ester (0.0924mmol) in 2.0mL ethyl acetate and 10mgs of 10% palladium on carbon was stirred under an atmosphere of hydrogen for 3 hours. The mixture was filtered, washing the solids with ethyl acetate, and concentrated *in vacuo* to yield 45mgs (99%) 1-methyl-4-(2-carboxyethyl)-7-(N-boc-5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione benzyl ester. <sup>1</sup>H NMR (CDCl<sub>3</sub>, δTMS) 7.63 (1H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, Ar-H *o*-CON), 7.35-7.27 (6H, Ph, *p*-CON), 7.06 (1H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, *m*-CON), 5.11 (2H, s, OCH<sub>2</sub>), 4.55 (1H, bs, BocNH), 3.98 (1H, d, <sup>2</sup>J<sub>HH</sub>=15Hz, NCHHCO), 3.94 (2H, q, <sup>3</sup>J<sub>HH</sub>=7Hz, NCH<sub>2</sub>), 3.80 (1H, d, <sup>2</sup>J<sub>HH</sub>=15Hz, NCHHCO), 3.35 (3H, s, NCH<sub>3</sub>), 3.08 (2H, q, <sup>3</sup>J<sub>HH</sub>=6.5Hz, BocNHCH<sub>2</sub>), 2.88-2.67 (2H, m, CH<sub>2</sub>CO<sub>2</sub>), 2.62 (2H, t, <sup>3</sup>J<sub>HH</sub>=8Hz, ArCH<sub>2</sub>), 1.65 (2H, p, <sup>3</sup>J<sub>HH</sub>=8Hz), 1.49 (2H, p, <sup>3</sup>J<sub>HH</sub>=8Hz), 1.42 (9H, s, *t*-Bu), 1.35 (2H, p, <sup>3</sup>J<sub>HH</sub>=8Hz). <sup>13</sup>C NMR (CDCl<sub>3</sub>, δTMS) 171.2, 168.9, 167.3, 140.1, 138.8, 135.6, 132.2, 130.2, 128.5, 128.4, 128.3, 120.9, 66.6, 52.3, 45.1, 40.4, 34.9, 34.8, 32.9, 30.7, 29.9, 28.4, 26.4.

Using the above procedure, but substituting the appropriate 7-alkynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione for 1-methyl-4-(2-carboxyethyl)-7-(N-boc-5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione benzyl ester there may be prepared, for example, the following compounds:

- 1-methyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 1-methyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-methyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-methyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-methyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-methyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- 1-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 (±)-1-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyglycynyl-4-benzyloxy)succinyl)]-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 (±)-1-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-benzyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 1-benzyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyglycynyl-4-benzyloxy)succinyl)]-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 (±)-1-benzyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 1-isopropyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-isopropyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 (±)-1-isopropyl-4-[2-(1-benzyloxyglycynyl-4-benzyloxy)succinyl)]-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 1-methyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- 1-methyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 5 (±)-1-methyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyl]-
- 10 3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 15 (±)-1-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyl]-3,4-
- 20 dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-benzyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 25 1-benzyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyl]-3,4-
- 30 dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 35 1-isopropyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-isopropyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-isopropyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 (±)-1-isopropyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-methyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 1-methyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 (±)-1-methyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 1-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 (±)-1-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 (±)-1-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-benzyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 1-benzyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-benzyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 (±)-1-benzyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-isopropyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-isopropyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 (±)-1-isopropyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 (±)-1-isopropyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-methyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 1-methyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 (±)-1-methyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 1-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 (±)-1-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 1-benzyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-benzyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-  
10 1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyglycinyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 (±)-1-benzyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-isopropyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-isopropyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-  
20 1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-benzyloxyglycinyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 (±)-1-isopropyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-methyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-  
30 piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-methyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 (±)-1-methyl-4-[2-(1-benzyloxyglycinyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-methyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 5 1-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 10 (±)-1-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 15 1-benzyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-benzyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 20 (±)-1-benzyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 25 (±)-1-benzyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-isopropyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-isopropyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 30 (±)-1-isopropyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 35 (±)-1-isopropyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- 1-methyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-methyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 (±)-1-methyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 (±)-1-methyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 1-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 (±)-1-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 1-benzyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-benzyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 (±)-1-benzyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 1-isopropyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,



- 1-isopropyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 5 (±)-1-isopropyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 10 1-methyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-methyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 15 (±)-1-methyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 20 (±)-1-methyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 25 1-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 30 (±)-1-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 35 1-benzyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-benzyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-benzyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 (±)-1-benzyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-isopropyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 1-isopropyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 (±)-1-isopropyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 1-methyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-methyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 (±)-1-methyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 (±)-1-methyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 1-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-phenyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 (±)-1-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-benzyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-benzyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 (±)-1-benzyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 (±)-1-benzyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-isopropyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 1-isopropyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 (±)-1-isopropyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 (±)-1-isopropyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-methyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-methyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 (±)-1-methyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-methyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 1-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 (±)-1-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 (±)-1-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-benzyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-benzyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 (±)-1-benzyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 (±)-1-benzyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-isopropyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-isopropyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 (±)-1-isopropyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-isopropyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 (±)-1-methyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl)]-7-[5-(N-Boc)-  
10 amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 (±)-1,3-diphenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 (±)-1,3-diphenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl)]-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 (±)-1,3-diphenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 (±)-1-benzyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl)]-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 (±)-1-isopropyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl)]-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyvaliny-4-benzyloxy)succinyl)]-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyphenylalany-4-benzyloxy)succinyl)]-7-[5-(N-Boc)-amino-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 (±)-1-methyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl)]-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 (±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyvaliny-4-benzyloxy)succinyl)]-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyphenylalany-4-benzyloxy)succinyl)]-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 (±)-1,3-diphenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 (±)-1,3-diphenyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl)]-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-benzyloxyvaliny-4-benzyloxy)succinyl)]-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 (±)-1,3-diphenyl-4-[2-(1-benzyloxyphenylalany-4-benzyloxy)succinyl)]-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 (±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[5-cyano-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 (±)-1-methyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 (±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 (±)-1,3-diphenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1,3-diphenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 (±)-1,3-diphenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 (±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 (±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[2-(3-cyanophenyl)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 (±)-1-methyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,



- (±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 (±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 (±)-1,3-diphenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 (±)-1,3-diphenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 (±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 (±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 5 (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[N-[2-(N-Boc)-aminoethane]-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 10 (±)-1-methyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-methyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-methyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 15 (±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 20 (±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1,3-diphenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1,3-diphenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 25 (±)-1,3-diphenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1,3-diphenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 30 (±)-1,3-diphenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1,3-diphenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 35 (±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-benzyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl)]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 5 (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 10 (±)-1-isopropyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl)]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 15 (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[3-keto-1-propyl-3-[4-(N-Boc)-1-piperazine]]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-[2-(N-
- 20 Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 25 (±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl)]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[2-[2-(N-
- 30 Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 35 (±)-1,3-diphenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl)]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1,3-diphenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 (±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 (±)-1-isopropyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[2-[2-(N-Boc)aminoethanethiol]-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 (±)-1-methyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 (±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 (±)-1,3-diphenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 (±)-1,3-diphenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 (±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 (±)-1-isopropyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]]-7-[2-(4-cyanothiophenol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-methyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
5 (±)-1-methyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
10 (±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
15 (±)-1,3-diphenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
20 (±)-1,3-diphenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1,3-diphenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
25 (±)-1,3-diphenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
30 (±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
(±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyglyciny-4-benzyloxy)succinyl]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
35 (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 5 (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 10 (±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-isopropyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 15 (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyglycinyl-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 20 (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[4-[2-(N-Boc)amino-6-methylpyrimidine-4-yl]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-methyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 25 (±)-1-methyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-methyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 30 (±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyglycinyl-4-benzyloxy)succinyl)]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl)]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 35 (±)-1-methyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl)]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1,3-diphenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1,3-diphenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- (±)-1,3-diphenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl)]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,

- (±)-1,3-diphenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 (±)-1,3-diphenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 5 (±)-1,3-diphenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 (±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 (±)-1-benzyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 10 (±)-1-benzyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 15 (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 (±)-1-benzyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 (±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-methyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 20 (±)-1-isopropyl-3-phenyl-4-(3-ethoxy-1-phenyl-3-oxo-1-propyl)-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 (±)-1-isopropyl-3-phenyl-4-[2-(1-*t*-butoxy-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 25 (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyglycyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyvalinyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 (±)-1-isopropyl-3-phenyl-4-[2-(1-benzyloxyphenylalanyl-4-benzyloxy)succinyl]-7-[5-[2-(N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 30 (N-Boc)aminopyridine-3-yl]-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
 b) 1-methyl-4-(2-carboxyethyl)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate was prepared using the method described in part  
 35 (g) of example 1. Thus, 45mgs 1-methyl-4-(2-carboxyethyl)-7-(N-boc-5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione benzyl ester yielded 36mgs (93%) 1-methyl-4-(2-carboxyethyl)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate (1/2" C-18 reverse-phase column, eluting with a solvent gradient of 10:90 acetonitrile(0.1% trifluoroacetic acid)/ water (0.1% trifluoroacetic acid), time 0 to 10 min, to



50:50 acetonitrile(0.1% trifluoroacetic acid)/water (0.1% trifluoroacetic acid), time 10 min to 40 min, flow=10ml/min,  $R_t=30.3$  min, un detection 254nm). HRMS (FAB) molecular ion  $m/z=348.1931$  ( cald.  $C_{18}H_{25}N_3O_4$ , 348.1923)  $^1H$  NMR ( $CDCl_3$ , dTMS) 7.39 (1H, bs, Ar-H o-CON), 7.34 (1H, d,  $^3J_{HH}=8$ Hz, p-CON), 7.16 (1H, d,  $^3J_{HH}=8$ Hz, m-CON), 4.00-3.90 (2H, m, NCHHCO, NCHHCH<sub>2</sub>), 3.67 (1H, d,  $^2J_{HH}=16$ Hz, NCHHCO), 3.56 (1H, dt,  $^2J_{HH}=14$ Hz,  $^3J_{HH}=6$ Hz, NCHHCH<sub>2</sub>) 3.17 (3H, s, NCH<sub>3</sub>), 2.82 (2H, t,  $^3J_{HH}=7$ Hz, H<sub>3</sub>NCH<sub>2</sub>), 2.67-2.47 (4H, m, CH<sub>2</sub>CO<sub>2</sub>, ArCH<sub>2</sub>), 1.57-1.42 (4H, bs), 1.16 (2H, p,  $^3J_{HH}=7$ Hz).

Using the above procedure, but substituting the appropriate 3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione for 1-methyl-4-(2-carboxyethyl)-7-(N-boc-5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione benzyl ester there could be prepared, for example, the following compounds:

- 1-methyl-4-(2-carboxy-1-methylethyl)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-methyl-4-(2-carboxy-1-phenylethyl)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-4-[succin-2-yl-(1-glycine)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-4-[succin-2-yl-(1-valine)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-phenyl-4-(2-carboxy-1-methylethyl)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-phenyl-4-(2-carboxy-1-phenylethyl)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-phenyl-4-[succin-2-yl-(1-glycine)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-phenyl-4-[succin-2-yl-(1-valine)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-benzyl-4-(2-carboxy-1-methylethyl)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-benzyl-4-(2-carboxy-1-phenylethyl)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1-benzyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-glycine)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
5 (±)-1-benzyl-4-[succin-2-yl-(1-valine)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-isopropyl-4-(2-carboxy-1-methylethyl)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
10 1-isopropyl-4-(2-carboxy-1-phenylethyl)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
15 (±)-1-isopropyl-4-[succin-2-yl-(1-glycine)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-valine)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
20 1-methyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-methyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
25 (±)-1-methyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
30 (±)-1-methyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-phenyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
35 1-phenyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1-phenyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-phenyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-benzyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-benzyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 (±)-1-benzyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1-benzyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-isopropyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 1-isopropyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1-isopropyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-aminoethane)-3-keto-1-propyn-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 1-methyl-4-(2-carboxy-1-methylethyl)-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-methyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-methyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-glycine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1-methyl-4-[succin-2-yl-(1-valine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
5 1-phenyl-4-(2-carboxy-1-methylethyl)-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-phenyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
10 (±)-1-phenyl-4-[succin-2-yl-(1-glycine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-valine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
15 (±)-1-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-benzyl-4-(2-carboxy-1-methylethyl)-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-benzyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
20 1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-glycine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
25 (±)-1-benzyl-4-[succin-2-yl-(1-valine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-isopropyl-4-(2-carboxy-1-methylethyl)-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
30 1-isopropyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
35 (±)-1-isopropyl-4-[succin-2-yl-(1-glycine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-valine)]-7-[3-keto-1-propyn-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1-isopropyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyn-3-(1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-methyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 1-methyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-
- 10 1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 1-phenyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-phenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-
- 20 1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-benzyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-benzyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-
- 30 1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-benzyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- 1-isopropyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-isopropyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
5 (±)-1-isopropyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
10 (±)-1-isopropyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-methyl-4-(2-carboxy-1-methylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
15 1-methyl-4-(2-carboxy-1-phenylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-glycine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
20 (±)-1-methyl-4-[succin-2-yl-(1-valine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-phenylalanine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
25 1-phenyl-4-(2-carboxy-1-methylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-phenyl-4-(2-carboxy-1-phenylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
30 (±)-1-phenyl-4-[succin-2-yl-(1-glycine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-valine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
35 (±)-1-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-benzyl-4-(2-carboxy-1-methylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- 1-benzyl-4-(2-carboxy-1-phenylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
5 (±)-1-benzyl-4-[succin-2-yl-(1-glycine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-valine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-phenylalanine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
10 1-isopropyl-4-(2-carboxy-1-methylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-isopropyl-4-(2-carboxy-1-phenylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
15 (±)-1-isopropyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-glycine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)]-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-valine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
20 (±)-1-isopropyl-4-[succin-2-yl-(1-phenylalanine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-methyl-4-(2-carboxy-1-methylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
25 1-methyl-4-(2-carboxy-1-phenylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-glycine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
30 (±)-1-methyl-4-[succin-2-yl-(1-valine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-phenylalanine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
35 1-phenyl-4-(2-carboxy-1-methylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-phenyl-4-(2-carboxy-1-phenylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-phenyl-4-[succin-2-yl-(1-glycine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1-phenyl-4-[succin-2-yl-(1-valine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-benzyl-4-(2-carboxy-1-methylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 1-benzyl-4-(2-carboxy-1-phenylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1-benzyl-4-[succin-2-yl-(1-glycine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-4-[succin-2-yl-(1-valine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-4-[succin-2-yl-(1-phenylalanine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 1-isopropyl-4-(2-carboxy-1-methylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-isopropyl-4-(2-carboxy-1-phenylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1-isopropyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-4-[succin-2-yl-(1-glycine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-4-[succin-2-yl-(1-valine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 (±)-1-isopropyl-4-[succin-2-yl-(1-phenylalanine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-methyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,



- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-(2-carboxy-1-methylethyl)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-(2-carboxy-1-phenylethyl)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 (±)-1,3-diphenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-glycine)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1,3-diphenyl-4-[succin-2-yl-(1-valine)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
5 (±)-1-methyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[N-(2-aminoethane)-3-keto-1-propyl-10 3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-aminoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-aminoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
15 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-aminoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
20 (±)-1,3-diphenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[N-(2-aminoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-aminoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-aminoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
25 (±)-1,3-diphenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-aminoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
30 (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[N-(2-aminoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-aminoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
35 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-aminoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-aminoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-aminoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[N-(2-aminoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-aminoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-aminoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-aminoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1-methyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[3-keto-1-propyl-3-(1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyl-3-(1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyl-3-(1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[3-keto-1-propyl-3-(1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[3-keto-1-propyl-3-(1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyl-3-(1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-(2-carboxy-1-methylethyl)-7-[3-keto-1-propyl-3-(1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyl-3-(1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 (±)-1,3-diphenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyl-3-(1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-glycine)]-7-[3-keto-1-propyl-3-(1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1,3-diphenyl-4-[succin-2-yl-(1-valine)]-7-[3-keto-1-propyl-3-(1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyl-3-(1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[3-keto-1-propyl-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyl-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyl-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[3-keto-1-propyl-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[3-keto-1-propyl-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyl-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[3-keto-1-propyl-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyl-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyl-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[3-keto-1-propyl-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[3-keto-1-propyl-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyl-3-(1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1-methyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1,3-diphenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
5 (±)-1,3-diphenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(2-aminoethanethiol)-1-ethyl]-  
10 3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
15 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
20 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
25 (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-aminoethanethiol)-1-ethyl]-  
30 3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(2-aminoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
35 (±)-1-methyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 (±)-1,3-diphenyl-4-(2-carboxy-1-methylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-(2-carboxy-1-phenylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1,3-diphenyl-4-[succin-2-yl-(1-glycine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-valine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 (±)-1,3-diphenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
5 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[4-(2-amino-6-methylpyrimidine-4-yl)-4-oxa-1-butyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
10 (±)-1-methyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
15 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
20 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-(2-carboxy-1-methylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-(2-carboxy-1-phenylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
25 (±)-1,3-diphenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-[succin-2-yl-(1-glycine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
30 (±)-1,3-diphenyl-4-[succin-2-yl-(1-valine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1,3-diphenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
35 (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 5 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 10 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 15 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[5-(2-aminopyridine-3-yl)-5-oxa-1-pentyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 c) 1-methyl-4-(2-carboxyethyl)-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione was prepared by the method described in Example 2.  
 Thus, 18mgs 1-methyl-4-(2-carboxyethyl)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione yielded 17mgs (87%) 1-methyl-4-(2-carboxyethyl)-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate (1/2" C-  
 25 18 reverse-phase column, eluting with a solvent gradient of 10:90 methanol(0.1% trifluoroacetic acid)/ water (0.1% trifluoroacetic acid), time 0 to 10 min, to 50:50 methanol(0.1% trifluoroacetic acid)/water (0.1% trifluoroacetic acid), time 10 min to 40 min, flow=10ml/min,  $R_t$ =38.5 min, un detection 254nm), HRMS (FAB) molecular ion  $m/z$ =390.2132 ( cald.  $C_{19}H_{27}N_5O_4$ , 390.2182)
- 30 Using the above procedure, but substituting the appropriate amino acid for 1-methyl-4-(2-carboxyethyl)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione there may be prepared, for example, the following compounds:  
 1-methyl-4-(2-carboxy-1-methylethyl)-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 35 1-methyl-4-(2-carboxy-1-phenylethyl)-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 (±)-1-methyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,



- (±)-1-methyl-4-[succin-2-yl-(1-glycine)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-4-[succin-2-yl-(1-valine)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1-methyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-phenyl-4-(2-carboxy-1-methylethyl)-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-phenyl-4-(2-carboxy-1-phenylethyl)-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 (±)-1-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-phenyl-4-[succin-2-yl-(1-glycine)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1-phenyl-4-[succin-2-yl-(1-valine)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-benzyl-4-(2-carboxy-1-methylethyl)-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 1-benzyl-4-(2-carboxy-1-phenylethyl)-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1-benzyl-4-[succin-2-yl-(1-glycine)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-4-[succin-2-yl-(1-valine)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 1-isopropyl-4-(2-carboxy-1-methylethyl)-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-isopropyl-4-(2-carboxy-1-phenylethyl)-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-isopropyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-4-[succin-2-yl-(1-glycine)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1-isopropyl-4-[succin-2-yl-(1-valine)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 1-methyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-methyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 (±)-1-methyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1-methyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-phenyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-phenyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 (±)-1-phenyl-4-[succin-2-yl-(1-*f*-butoxy)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1-phenyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 1-benzyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-benzyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-benzyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1-benzyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-isopropyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
5 1-isopropyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
10 (±)-1-isopropyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
15 1-methyl-4-(2-carboxy-1-methylethyl)-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-methyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
20 (±)-1-methyl-4-[succin-2-yl-(1-glycine)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-valine)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
25 (±)-1-methyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-phenyl-4-(2-carboxy-1-methylethyl)-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-phenyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
30 (±)-1-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-glycine)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-valine)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
35 (±)-1-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- 1-benzyl-4-(2-carboxy-1-methylethyl)-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-benzyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1-benzyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-4-[succin-2-yl-(1-glycine)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-4-[succin-2-yl-(1-valine)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 (±)-1-benzyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-isopropyl-4-(2-carboxy-1-methylethyl)-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 1-isopropyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-4-[succin-2-yl-(1-glycine)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 (±)-1-isopropyl-4-[succin-2-yl-(1-valine)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 1-methyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-methyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 (±)-1-methyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-methyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-phenyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- 1-phenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1-phenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-
- 10 dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-benzyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-benzyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1-benzyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-
- 20 dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-isopropyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 1-isopropyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-
- 30 dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-
- 35 3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-(2-carboxy-1-methylethyl)-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 (±)-1,3-diphenyl-4-(2-carboxy-1-phenylethyl)-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1,3-diphenyl-4-[succin-2-yl-(1-glycine)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-valine)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 (±)-1,3-diphenyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1,3-diphenyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

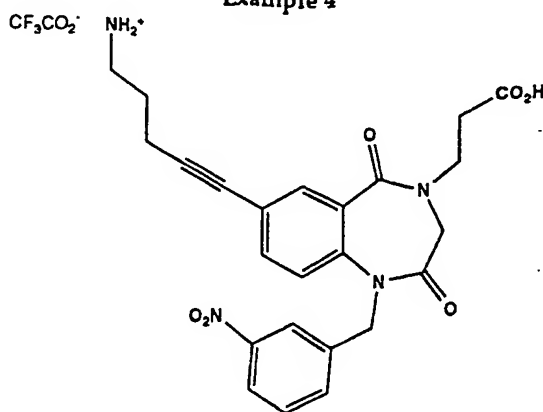
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[N-(2-guanidinoethane)-3-keto-1-propyl-3-amine]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[3-keto-1-propyl-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyl-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-(2-carboxy-1-methylethyl)-7-[3-keto-1-propyl-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 (±)-1,3-diphenyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyl-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-glycine)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1,3-diphenyl-4-[succin-2-yl-(1-valine)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperazine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,



- (±)-1,3-diphenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[3-keto-1-propyl-3-(4-amidino-1-piperizine)]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1,3-diphenyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1,3-diphenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 (±)-1,3-diphenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(2-guanidinoethanethiol)-1-ethyl]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate.

## Example 4



1-(3-nitrobenzyl)-4-(2-carboxyethyl)-7-(5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate.

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a) N-(2-amino-5-iodobenzoyl)-b-alanine ethyl ester was prepared using the method described in part (b) of Example 1. Thus, from 21.0 grams of 6-iodoisatoic anhydride (0.73 mol) and 11.3 grams of b-alanine ethyl ester hydrogenschloride (0.73 mol) was prepared 9.1 grams (35%) N-(2-amino-5-iodobenzoyl)-b-alanine ethyl ester (mp=83-85°C, TLC, SiO<sub>2</sub>, 1:1 EtOAc/hexane, R<sub>f</sub>=0.46, un positive). <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 7.57 (1H, d, <sup>4</sup>J<sub>HH</sub> = 2Hz, Ar-H *o*-CON), 7.42 (1H, dd, <sup>3</sup>J<sub>HH</sub>=9Hz, <sup>4</sup>J<sub>HH</sub>=2Hz, Ar-H *p*-CON), 6.63 (1H, bs, CONH), 6.44 (1H, d, <sup>3</sup>J<sub>HH</sub>=9Hz, Ar-H *m*-CON), 5.57 (2H, bs, NH<sub>2</sub>), 4.10 (2H, q, <sup>3</sup>J<sub>HH</sub>=7Hz, OCH<sub>2</sub>), 3.63 (2H, q, <sup>3</sup>J<sub>HH</sub>=6Hz, NCH<sub>2</sub>), 2.63 (2H, t, <sup>3</sup>J<sub>HH</sub>=6Hz, CH<sub>2</sub>CO<sub>2</sub>), 1.14 (3H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, CH<sub>3</sub>).

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b) A magnetically stirred solution of 1.2 grams N-(2-amino-5-iodobenzoyl)-b-alanine ethyl ester (3.35 mmol), 0.43mL 2,6-lutidine (3.7 mmol), 0.713 grams 3-nitrobenzyl bromide (3.3 mmol), and 20mL dimethylformamide, under an atmosphere of nitrogen, was heated to 50°C for 24 hours. The reaction mixture was allowed to cool to room temperature, poured over 100mL 10% citric acid, and extracted 3X75mL ether. The combined organics were dried over magnesium sulfate, filtered and concentrated *in vacuo*. The resulting residue was purified by column chromatography, using silica gel, eluting with a solvent gradient of 1:4 ethyl acetate/hexane to 1:1 ethyl acetate/hexane isolating the yellow band (TLC, SiO<sub>2</sub>, 1:1 ethyl acetate/hexane, R<sub>f</sub>=0.66, un positive) characterized as N-[2-(3-nitrobenzyl)-5-iodobenzoyl]-b-alanine ethyl ester (1.1 grams, 66%, mp=104-105°C).

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c) To a magnetically stirred solution of 1.0 grams N-[2-(3-nitrobenzyl)-5-iodobenzoyl]-b-alanine ethyl ester (2.0 mmol), 5mL methylene chloride, 0.56mL triethylamine (4.0 mmol) at -30°C under an atmosphere of nitrogen was slowly added, *via* an addition funnel, 0.26mL a-bromoacetyl bromide (3.0 mmol) as a solution in 4mL methylene chloride. After 2 hours, the mixture was diluted with 35mL methylene chloride and washed 2X75mL 10% citric acid,

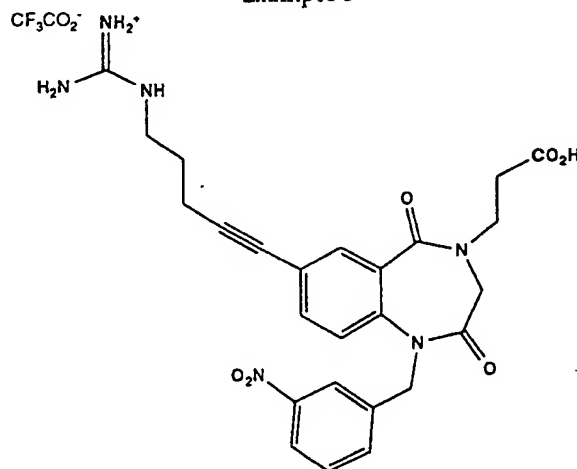
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2X75mL sat. sodium bicarbonate, dried over magnesium sulfate, filtered and concentrated *in vacuo*. The resulting residue was dissolved in 5mL dimethylformamide and added, *via* an addition funnel, to a slurry of 72mgs sodium hydride (3.0 mmol) in 1.0mL dimethylformamide cooled to 0°C. The reaction was allowed to warm to room temperature. After 2 hours, the mixture was poured over 100mL of an ice cooled solution of 10% citric acid and extracted 3X75mL ethyl acetate. The combined organics were washed with 1X50mL sat. sodium bicarbonate, dried over magnesium sulfate, and concentrated *in vacuo*. The resulting residue was further purified by column chromatography, using silica gel, eluting with a solvent gradient of 40:60 ethyl acetate/hexane to 70:30 ethyl acetate/hexane to yield 0.45 grams (55%) 1-(3-nitrobenzyl)-4-(2-carboxyethyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester (TLC, SiO<sub>2</sub>, 1:1 ethyl acetate/hexane, R<sub>f</sub>=0.39, un positive) and 0.18 grams of recovered N-[2-(3-nitrobenzyl)-5-iodobenzoyl]-b-alanine ethyl ester. <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 8.09 (1H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, o-CON), 8.06 (1H, dt, <sup>3</sup>J<sub>HH</sub>=7Hz, <sup>4</sup>J<sub>HH</sub>=2Hz, NCH<sub>2</sub>(C4-Ar-H)), 8.00 (1H, bs, NCH<sub>2</sub>(C2-Ar-H)), 7.73 (1H, dd, <sup>3</sup>J<sub>HH</sub>=9Hz, <sup>4</sup>J<sub>HH</sub>=2Hz, Ar-H p-CON), 7.5-7.4 [2H, m, NCH<sub>2</sub>(C5,C6-Ar-H)], 6.88 (1H, d, <sup>3</sup>J<sub>HH</sub>=9Hz, Ar-H m-CON), 5.20 (1H, d, <sup>2</sup>J<sub>HH</sub>=16Hz, NCHHAr), 5.03 (1H, d, <sup>2</sup>J<sub>HH</sub>=16Hz, NCHHAr), 4.15 (1H, d, <sup>2</sup>J<sub>HH</sub>=15Hz, NCHHCO), 4.10 (2H, q, <sup>3</sup>J<sub>HH</sub>=7Hz, OCH<sub>2</sub>), 3.96 (1H, d, <sup>2</sup>J<sub>HH</sub>=15Hz, NCHHCO), 3.98-3.92 (2H, m, NCH<sub>2</sub>CH<sub>2</sub>), 2.71 (1H, dt, <sup>2</sup>J<sub>HH</sub>=16Hz, <sup>3</sup>J<sub>HH</sub>=8Hz, CHHCO<sub>2</sub>), 2.63 (1H, dt, <sup>2</sup>J<sub>HH</sub>=16Hz, <sup>3</sup>J<sub>HH</sub>=8Hz, CHHCO<sub>2</sub>), 1.22 (3H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, CH<sub>3</sub>). <sup>13</sup>C NMR (CDCl<sub>3</sub>, dTMS) 171.2, 168.3, 165.3, 148.4, 141.0, 139.7, 139.0, 138.5, 132.8, 131.1, 129.9, 122.9, 122.8, 121.9, 90.5, 60.8, 52.1, 50.3, 45.2, 32.7, 14.1.

d) 1-(3-nitrobenzyl)-4-(2-carboxyethyl)-7-(N-boc-5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester was prepared using the method described in part (f) of Example 1. Thus, from 0.205 grams of 1-(3-nitrobenzyl)-4-(2-carboxyethyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester was prepared 168.2mgs (57%) 1-(3-nitrobenzyl)-4-(2-carboxyethyl)-7-(N-boc-5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester (TLC, SiO<sub>2</sub>, 1:1 ethyl acetate/hexane, R<sub>f</sub>=0.36, un positive). <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 8.09 [1H, dt, <sup>3</sup>J<sub>HH</sub>=7Hz, <sup>4</sup>J<sub>HH</sub>=2Hz, NCH<sub>2</sub>(C4-Ar-H)], 8.02 [1H, bs, NCH<sub>2</sub>(C2-Ar-H)], 7.81 (1H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, o-CON), 7.50-7.35 [3H, m, NCH<sub>2</sub>(C5,C6-Ar-H), Ar-H p-CON], 7.04 (1H, d, <sup>3</sup>J<sub>HH</sub>=9Hz, Ar-H m-CON), 5.22 (1H, d, <sup>2</sup>J<sub>HH</sub>=16Hz, NCHHAr), 5.04 (1H, d, <sup>2</sup>J<sub>HH</sub>=16Hz, NCHHAr), 4.70 (1H, bs, BocNH), 4.40 (1H, d, <sup>2</sup>J<sub>HH</sub>=12Hz, NCHHCO), 4.12 (2H, q, <sup>3</sup>J<sub>HH</sub>=7Hz, OCH<sub>2</sub>), 3.96 (1H, d, <sup>2</sup>J<sub>HH</sub>=12Hz, NCHHCO), 4.0-3.8 (2H, m, NCH<sub>2</sub>CH<sub>2</sub>), 3.23 (2H, q, <sup>3</sup>J<sub>HH</sub>=6Hz, BocNHCH<sub>2</sub>), 2.72 (1H, dt, <sup>2</sup>J<sub>HH</sub>=17Hz, <sup>3</sup>J<sub>HH</sub>=8Hz, CHHCO<sub>2</sub>), 2.62 (1H, dt, <sup>2</sup>J<sub>HH</sub>=17Hz, <sup>3</sup>J<sub>HH</sub>=8Hz, CHHCO<sub>2</sub>), 2.41 (2H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, CJCCH<sub>2</sub>), 1.74 (2H, p, <sup>3</sup>J<sub>HH</sub>=7Hz, CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>), 1.40 (9H, s, t-Bu), 1.24 (3H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, CH<sub>3</sub>). <sup>13</sup>C NMR (CDCl<sub>3</sub>, dTMS) 171.2, 168.4, 166.2, 155.9, 148.4, 138.6, 138.2, 134.9, 134.2, 132.8, 129.9, 129.5, 122.7, 122.4, 122.0, 121.2, 91.4, 79.2, 60.8, 52.2, 50.4, 45.2, 39.8, 32.7, 28.7, 28.4, 16.8, 14.1.

- e) 1-(3-nitrobenzyl)-4-(2-carboxyethyl)-7-(5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate was prepared using the method described in part (g) of example 1. Thus, 84 mgs 1-(3-nitrobenzyl)-4-(2-carboxyethyl)-7-(N-boc-5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester (0.142 mmol) yielded 75mgs (91%) 1-(3-nitrobenzyl)-4-(2-carboxyethyl)-7-(5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate (1/2" C-18 reverse-phase column, eluting with a solvent gradient of 10:90 methanol(0.1% trifluoroacetic acid)/ water (0.1% trifluoroacetic acid), time 0 to 10 min, to 50:50 methanol(0.1% trifluoroacetic acid)/water (0.1% trifluoroacetic acid), time 10 min to 40 min, flow=10ml/min,  $R_t$ =45.1 min, un detection 254nm). HRMS (FAB) molecular ion  $m/z$ =465.1744 ( calcd.  $C_{24}H_{25}N_4O_6$ , 465.1774)  $^1H$  NMR ( $CD_3$  CN, dTMS) 8.07 [1H, dt,  $^4J_{HH}$ =2Hz,  $^3J_{HH}$ =7Hz,  $NCH_2(C4-Ar-H)$ ], 8.02 [1H, bs,  $NCH_2(C2-Ar-H)$ ], 7.73 (1H, d,  $^4J_{HH}$ =2Hz,  $o-CON$ ), 7.60-7.50 [3H, m,  $NCH_2(C5,C6-Ar-H)$ ,  $Ar-H$   $p-CON$ ], 7.25 (1H, d,  $^3J_{HH}$ =9Hz,  $Ar-H$   $m-CON$ ), 5.40 (1H, d,  $^2J_{HH}$ =15Hz,  $NCHHAr$ ), 5.00 (1H, d,  $^2J_{HH}$ =15Hz,  $NCHHAr$ ), 4.13 (1H, d,  $^2J_{HH}$ =12Hz,  $NCHHCO$ ), 4.0-3.6 (3H, m,  $NCHHCO$ ,  $NCH_2CH_2$ ), 3.1 (2H, bs,  $NH_3CH_2$ ), 2.65 (2H, t,  $^3J_{HH}$ =7Hz,  $CH_2CO_2$ ), 2.41 (2H, t,  $^3J_{HH}$ =7Hz,  $C(CCH_2)$ ), 1.9 (2H, p,  $^3J_{HH}$ =7Hz,  $CH_2CH_2CH_2$ ).

Example 5



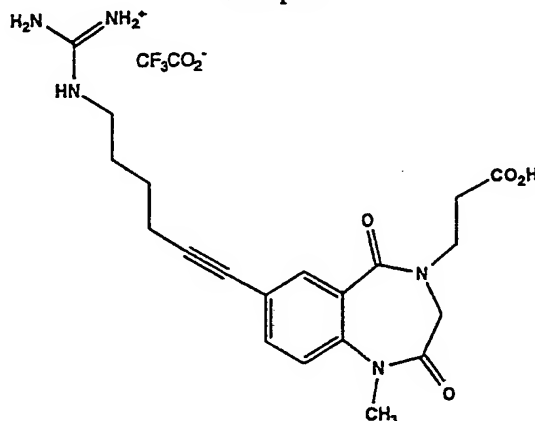
20

**1-(3-nitrobenzyl)-4-(2-carboxyethyl)-7-(5-guanidino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate.**

- 1-(3-nitrobenzyl)-4-(2-carboxyethyl)-7-(5-guanidino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate was prepared using the method described in Example 2. Thus, 30mgs of 1-(3-nitrobenzyl)-4-(2-carboxyethyl)-7-(5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate yielded 14mgs (44%) of 1-(3-nitrobenzyl)-4-(2-carboxyethyl)-7-(5-guanidino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate (1/2" C-18 reverse-phase column, eluting with a solvent gradient of 20:80 methanol(0.1% trifluoroacetic acid)/ water (0.1% trifluoroacetic

acid), time 0 to 10 min, to 70:30 methanol(0.1% trifluoroacetic acid)/water (0.1% trifluoroacetic acid), time 10 min to 40 min, flow=10ml/min,  $R_t=33.8$  min, un detection 254nm). HRMS (FAB) molecular ion  $m/z=507.2007$  ( cald.  $C_{25}H_{27}N_6O_6$ , 507.1992)  $^1H$  NMR ( $CD_3CN/CD_3OD$ , dTMS) 8.04 [1H, d,  $^3J_{HH}=7Hz$ ,  $NCH_2(C4-Ar-H)$ ], 7.98 [1H, bs,  $NCH_2(C2-Ar-H)$ ], 7.68 (1H, d,  $^4J_{HH}=2Hz$ ,  $o-CON$ ), 7.60-7.40 [3H, m,  $NCH_2(C5,C6-Ar-H)$ ,  $Ar-H p-CON$ ], 7.25 (1H, d,  $^3J_{HH}=9Hz$ ,  $Ar-H m-CON$ ), 5.37(1H, d,  $^2J_{HH}=15Hz$ ,  $NCHHAr$ ), 5.00 (1H, d,  $^2J_{HH}=15Hz$ ,  $NCHHAr$ ), 4.13 (1H, d,  $^2J_{HH}=12Hz$ ,  $NCHHCO$ ), 4.0-3.6 (3H, m,  $NCHHCO$ ,  $NCH_2CH_2$ ), 3.24 (2H, bs,  $H_2N(H_2N=)NHCH_2$ ), 2.65 (2H, t,  $CH_2CO_2$ ), 2.41 (2H, t,  $C[CCH_2]$ ), 1.9 (2H, p,  $^3J_{HH}=7Hz$ ,  $CH_2CH_2CH_2$ ).

Example 6



*1-(methyl)-4-(2-carboxyethyl)-7-(6-guanidino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate.*

- 15 a) N-(2-methylamino-5-iodobenzoyl)-b-alanine ethyl ester was prepared using the method described in part (c) of Example 1. Thus, 0.5grams N-(2-amino-5-iodobenzoyl)-b-alanine ethyl ester yielded 204.5 mgs (39%) of N-(2-methylamino-5-iodobenzoyl)-b-alanine ethyl ester (TLC,  $SiO_2$ , 1:1 ethyl acetate/hexane  $R_f=0.67$ ) and 253.2mgs of a 7:1 mixture of N-(2-amino-5-iodobenzoyl)-b-alanine ethyl ester and N-(2-dimethylamino-5-iodobenzoyl)-b-alanine ethyl ester (TLC,  $SiO_2$ , 1:1 ethyl acetate/hexane  $R_f=0.46$ , un positive).  $^1H$  NMR ( $CDCl_3$ , dTMS) 7.6-7.5 (2H, m,  $Ar-H o,p-CON$ ), 7.43 (1H, bs,  $NHMe$ ), 6.62 (1H, bt,  $CONH$ ), 6.42 (1H, d,  $^3J_{HH}=9Hz$ ,  $Ar-H m-CON$ ), 4.18 (2H, q,  $^3J_{HH}=7Hz$ ,  $OCH_2$ ), 3.63 (2H, q,  $^3J_{HH}=6Hz$ ,  $NCH_2$ ), 2.81 (3H, d,  $^3J_{HH}=5Hz$ ,  $NCH_3$ ), 2.61 (2H, t,  $^3J_{HH}=6Hz$ ,  $CH_2CO_2$ ), 1.28 (3H, t,  $^3J_{HH}=7Hz$ ,  $CH_3$ ).

- 25 b) 1-methyl-4-(2-carboxyethyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester was prepared using the method described in part (d) of Example 1. Thus, 1.9 grams of N-(2-methylamino-5-iodobenzoyl)-b-alanine ethyl ester yielded 0.91 grams (43%) of 1-methyl-4-(2-carboxyethyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester (TLC,  $SiO_2$ , 1:1 ethyl acetate/hexane  $R_f=0.36$ , un positive).  $^1H$  NMR ( $CDCl_3$ , dTMS)
- 30

- 8.14 (1H, d,  $^4J_{HH}=2\text{Hz}$ , Ar-H *o*-CON), 7.79 (1H, dd,  $^4J_{HH}=2\text{Hz}$ ,  $^3J_{HH}=9\text{Hz}$ , Ar-H *p*-CON), 6.92 (1H, d,  $^3J_{HH}=9\text{Hz}$ , Ar-H *m*-CON), 4.13 (2H, q,  $^3J_{HH}=7\text{Hz}$ , OCH<sub>2</sub>), 4.01 (1H, d,  $^2J_{HH}=15\text{Hz}$ , COCHHN), 3.90 (2H, dt,  $^3J_{HH}=8\text{Hz}$ ,  $J_{HH}=2\text{Hz}$ , NCH<sub>2</sub>), 3.87 (1H, d,  $^2J_{HH}=15\text{Hz}$ , COCHHN), 3.34 (3H, s, NCH<sub>3</sub>), 2.74 (1H, dt,  $^2J_{HH}=17\text{Hz}$ ,  $^3J_{HH}=7\text{Hz}$ , CHHCO<sub>2</sub>), 2.63 (1H, dt,  $^2J_{HH}=17\text{Hz}$ ,  $^3J_{HH}=7\text{Hz}$ , CHHCO<sub>2</sub>), 1.24 (3H, t,  $^3J_{HH}=7\text{Hz}$ , CH<sub>3</sub>).

c) To a magnetically stirred slurry of 2.94 grams lithium aluminum hydride (0.077 mol) in 150mL ether cooled to 0°C was slowly added, *via* an addition funnel, 6 grams of 5-cyano-1-pentyne (0.065 mol) in 15mL ether. The reaction mixture was allowed to warm to room temperature. After 1 hour, the reaction was cooled to 0°C and quenched with 3.1 mL water, 2.33 mL 20% sodium hydroxide, 10.83mL water added in succession. After an additional 30 minutes, the slurry was filtered, washing the salts with 40mL ether. The ether was distilled at atmospheric pressure to leave an oil, which was further purified by Kugelrohr distillation (oven temp. ca. 145°C) to yield 3.99 grams 6-amino-1-hexyne (64%).

d) To a magnetically stirred solution of 3.99 grams of 6-amino-1-hexyne (0.0413 mol) in 50mL tetrahydrofuran was added 9.1 grams of di-*tert*-butyl dicarbonate (0.042 mol). After 2 hours, the reaction mixture was concentrated *in vacuo* and the resulting residue chromatographed, using silica gel, eluting with 1:1 ether/hexane to yield 3.4 grams (42%) N-boc-6-amino-1-hexyne (TLC, SiO<sub>2</sub>, 1:1 ether/hexane,  $R_f=0.5$ , ninhydrin char). <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 4.55 (1H, bs, BocNH), 3.13 (2H, bq,  $^3J_{HH}=5\text{Hz}$ , BocNHCH<sub>2</sub>), 2.21 (2H, dt,  $^4J_{HH}=3\text{Hz}$ ,  $^3J_{HH}=7\text{Hz}$ , CJCCH<sub>2</sub>), 1.95 (1H, t,  $^4J_{HH}=3\text{Hz}$ , CJCCH), 1.6 (4H, m, CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>), 1.42 (9H, s, *t*-Bu).

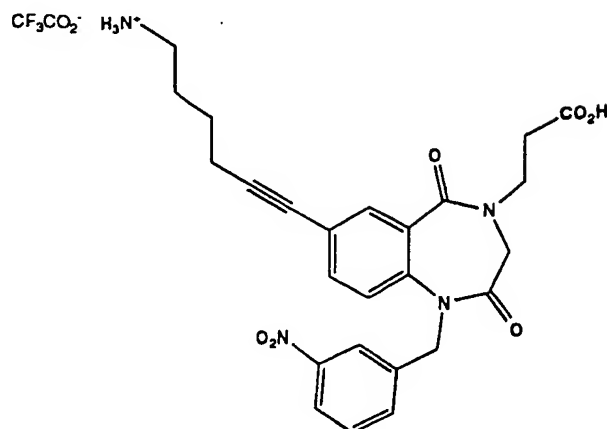
e) 1-(methyl)-4-(2-carboxyethyl)-7-(N-boc-6-amino-1-hexynyl)-3,4-dihydro-1*H*-1,4-benzodiazepine-2,5-dione ethyl ester was prepared using the method described in part (f) of Example 1. Thus, 0.126 grams of 1-methyl-4-(2-carboxyethyl)-7-iodo-3,4-dihydro-1*H*-1,4-benzodiazepine-2,5-dione ethyl ester and 0.119 grams of N-boc-6-amino-1-hexyne yielded 0.124 grams (84%) 1-(methyl)-4-(2-carboxyethyl)-7-(N-boc-6-amino-1-hexynyl)-3,4-dihydro-1*H*-1,4-benzodiazepine-2,5-dione ethyl ester (TLC, SiO<sub>2</sub>, 1:1 ethyl acetate/hexane,  $R_f=0.36$ , un positive). <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 7.84 (1H, d,  $^4J_{HH}=2\text{Hz}$ , Ar-H *o*-CON), 7.47 (1H, dd,  $^4J_{HH}=2\text{Hz}$ ,  $^3J_{HH}=9\text{Hz}$ , Ar-H *p*-CON), 7.07 (1H, d,  $^3J_{HH}=9\text{Hz}$ , Ar-H *m*-CON), 4.60 (1H, bs, NH), 4.12 (2H, q,  $^3J_{HH}=7\text{Hz}$ , OCH<sub>2</sub>), 4.0 (1H, d,  $^2J_{HH}=15\text{Hz}$ , NCHHCO), 3.90 (2H, t,  $^3J_{HH}=7\text{Hz}$ , NCH<sub>2</sub>), 3.84 (1H, d,  $^2J_{HH}=15\text{Hz}$ , NCHHCO), 3.26 (3H, s, NCH<sub>3</sub>), 3.15 (2H, q,  $^3J_{HH}=7\text{Hz}$ , BocNHCH<sub>2</sub>), 2.73 (1H, dt,  $^2J_{HH}=16\text{Hz}$ ,  $^3J_{HH}=9\text{Hz}$ , CHHCO<sub>2</sub>), 2.63 (1H, dt,  $^2J_{HH}=16\text{Hz}$ ,  $^3J_{HH}=9\text{Hz}$ , CHHCO<sub>2</sub>), 2.40 (2H, t,  $^3J_{HH}=6\text{Hz}$ , CJCCH<sub>2</sub>), 1.62 (4H, m, CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>), 1.24 (3H, t,  $^3J_{HH}=7\text{Hz}$ , CH<sub>3</sub>).

f) 1-(methyl)-4-(2-carboxyethyl)-7-(6-amino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate was prepared using the method described in part (g) of Example 1. Thus, 62mgs of 1-(methyl)-4-(2-carboxyethyl)-7-(N-boc-6-amino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester yielded 50mgs (83%) 1-(methyl)-4-(2-carboxyethyl)-7-(6-amino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate (1/2" C-18 reverse-phase column, eluting with a solvent gradient of 20:80 methanol(0.1% trifluoroacetic acid)/ water (0.1% trifluoroacetic acid), time 0 to 10 min, to 70:30 methanol(0.1% trifluoroacetic acid)/water (0.1% trifluoroacetic acid), time 10 min to 40 min, flow=10ml/min,  $R_t$ =26.2 min, un detection 254nm). HRMS (FAB) molecular ion  $m/z$ =358.1761 ( cald.  $C_{19}H_{24}N_3O_4$ , 358.1767).  $^1H$  NMR ( $D_2O$ ) 7.47 (1H, d,  $^4J_{HH}$ =2Hz, Ar-H *o*-CON), 7.37 (1H, dd,  $^4J_{HH}$ =2Hz,  $^3J_{HH}$ =9Hz, Ar-H *p*-CON), 7.12 (1H, d,  $^3J_{HH}$ =9Hz, Ar-H *m*-CON), 3.95-3.8 (2H, m, NCHHCO, NCHHCH<sub>2</sub>), 3.63 (1H, d,  $^2J_{HH}$ =15Hz, NCHHCO), 3.52 (1H, dt,  $^2J_{HH}$ =15Hz,  $^3J_{HH}$ =5Hz, NCHHCO<sub>2</sub>), 3.16 (3H, s, NCH<sub>3</sub>), 2.85 (2H, t,  $^3J_{HH}$ =7Hz, NH<sub>3</sub>CH<sub>2</sub>), 2.6-2.4 (2H, m,  $\dot{C}H_2CO_2$ ), 2.28 (2H, t,  $^3J_{HH}$ =6Hz, C(CCH<sub>2</sub>), 1.62 & 1.47 (each 2H, p,  $^3J_{HH}$ =6Hz, CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>).

g) 1-(methyl)-4-(2-carboxyethyl)-7-(6-guanidino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate was prepared by the method described in Example 2. Thus, from 25mgs 1-(methyl)-4-(2-carboxyethyl)-7-(6-amino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate was prepared 22mgs (71%) of 1-(methyl)-4-(2-carboxyethyl)-7-(6-guanidino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate (1/2" C-18 reverse-phase column, eluting with a solvent gradient of 20:80 methanol(0.1% trifluoroacetic acid)/ water (0.1% trifluoroacetic acid), time 0 to 10 min, to 70:30 methanol(0.1% trifluoroacetic acid)/water (0.1% trifluoroacetic acid), time 10 min to 40 min, flow=10ml/min,  $R_t$ =28.6 min, un detection 254nm). HRMS (FAB) molecular ion  $m/z$ =400.1956 ( cald.  $C_{20}H_{26}N_5O_4$ , 400.1985).  $^1H$  NMR ( $D_2O$ ) 7.58 (1H, d,  $^4J_{HH}$ =2Hz, Ar-H *o*-CON), 7.43 (1H, dd,  $^4J_{HH}$ =2Hz,  $^3J_{HH}$ =9Hz, Ar-H *p*-CON), 7.20 (1H, d,  $^3J_{HH}$ =9Hz, Ar-H *m*-CON), 4.05-3.95 (2H, m, NCHHCO<sub>2</sub>, NCHHCH<sub>2</sub>), 3.71 (1H, d,  $^2J_{HH}$ =15Hz, NCHHCO), 3.56 (1H, dt,  $^2J_{HH}$ =15Hz,  $^3J_{HH}$ =5Hz, NCHHCO<sub>2</sub>), 3.19 (3H, s, NCH<sub>3</sub>), 3.06 (2H, t,  $^3J_{HH}$ =7Hz, H<sub>2</sub>N(H<sub>2</sub>N=)CNHCH<sub>2</sub>), 2.7-2.5 (2H, m, CH<sub>2</sub>CO<sub>2</sub>), 2.32 (2H, t,  $^3J_{HH}$ =6Hz, C(CCH<sub>2</sub>), 1.60 & 1.49 (each 2H, p,  $^3J_{HH}$ =6Hz, CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>).



## Example 7



5            **1-(3-nitrobenzyl)-4-(2-carboxyethyl)-7-(6-amino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate.**

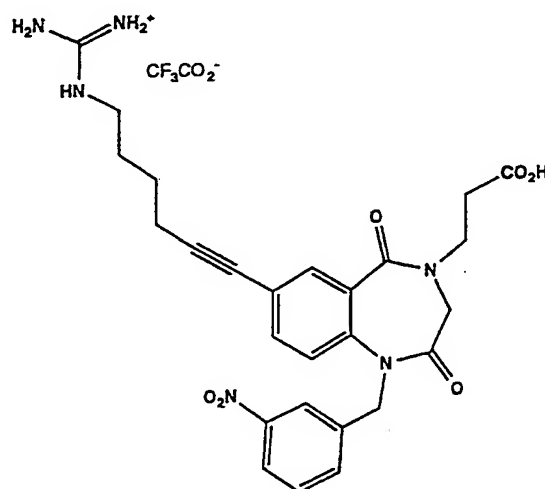
a)        1-(3-nitrobenzyl)-4-(2-carboxyethyl)-7-(N-boc-6-amino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester was prepared using the method described in part (f) of Example 1. Thus, from 0.26 grams of 1-(3-nitrobenzyl)-4-(2-carboxyethyl)-7-iodo-  
 10        3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester and 0.19 grams of N-boc-6-amino-1-hexyne was prepared 226.3mgs (78%) 1-(3-nitrobenzyl)-4-(2-carboxyethyl)-7-(N-boc-6-amino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester (TLC, SiO<sub>2</sub>, 1:1 ethyl acetate/hexane, R<sub>f</sub>=0.34, un positive). <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 8.09 [1H, dt, <sup>3</sup>J<sub>HH</sub>=7Hz, <sup>4</sup>J<sub>HH</sub>=2Hz, NCH<sub>2</sub>(C4-Ar-H)], 8.02 [1H, bs, NCH<sub>2</sub>(C2-Ar-H)], 7.83 (1H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, o-CON), 7.50-7.35 [3H, m, NCH<sub>2</sub>(C5,C6-Ar-H), Ar-H p-CON], 7.05 (1H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, Ar-H m-CON), 5.22 (1H, d, <sup>2</sup>J<sub>HH</sub>=16Hz, NCHHAr), 5.02 (1H, d, <sup>2</sup>J<sub>HH</sub>=16Hz, NCHHAr), 4.60 (1H, bs, BocNH), 4.20-4.08 (3H, m, NCHHCO, OCH<sub>2</sub>), 4.0-3.8 (3H, m, NCH<sub>2</sub>CH<sub>2</sub>, NCHHCO), 3.16 (2H, q, <sup>3</sup>J<sub>HH</sub>=6Hz, BocNHCH<sub>2</sub>), 2.72 (1H, dt, <sup>2</sup>J<sub>HH</sub>=16Hz, <sup>3</sup>J<sub>HH</sub>=8Hz, CHHCO<sub>2</sub>), 2.62 (1H, dt, <sup>2</sup>J<sub>HH</sub>=16Hz, <sup>3</sup>J<sub>HH</sub>=8Hz, CHHCO<sub>2</sub>), 2.39 (2H, t, <sup>3</sup>J<sub>HH</sub>=6Hz, C[CCH<sub>2</sub>]), 1.60 (4H, bs, CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>), 1.41 (9H, s, t-Bu), 1.24 (3H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, CH<sub>3</sub>).

b)        1-(3-nitrobenzyl)-4-(2-carboxyethyl)-7-(6-amino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate was prepared using the method described in part (g) of Example 1. Thus, from 0.11 grams of 1-(3-nitrobenzyl)-4-(2-carboxyethyl)-7-(N-boc-6-amino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester was prepared  
 25        90mgs (79%) of 1-(3-nitrobenzyl)-4-(2-carboxyethyl)-7-(6-amino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate (1/2" C-18 reverse-phase column, eluting with a solvent gradient of 20:80 methanol(0.1% trifluoroacetic acid)/ water (0.1% trifluoroacetic acid), time 0 to 10 min, to 70:30 methanol(0.1% trifluoroacetic acid)/water (0.1%  
 30        trifluoroacetic acid), time 0 to 10 min, to 70:30 methanol(0.1% trifluoroacetic acid)/water (0.1%

trifluoroacetic acid), time 10 min to 40 min, flow=10ml/min,  $R_t$ =39.2 min, un detection 254nm).  
 HRMS (FAB) molecular ion  $m/z$ =479.1911 ( cald.  $C_{25}H_{27}N_4O_6$ , 479.1931).  $^1H$  NMR ( $D_2O$ )  
 7.7 [1H, bd,  $^3J_{HH}$ =8Hz,  $NCH_2(C4-Ar-H)$ ], 7.58 [1H, bs,  $NCH_2(C2-Ar-H)$ ], 7.41 (1H, d,  
 $^4J_{HH}$ =2Hz,  $o$ -CON), 7.25-7.05 [3H, m,  $NCH_2(C5,C6-Ar-H)$ , Ar-H  $p$ -CON], 7.01 (1H, d,  
 5  $^3J_{HH}$ =9Hz, Ar-H  $m$ -CON), 5.09 (1H, d,  $^2J_{HH}$ =16Hz,  $NCHHAr$ ), 4.8 (1H, d,  $NCHHAr$ ,  
 overlapping with HOD), 4.00 (1H, d,  $^2J_{HH}$ =14Hz,  $NCHHCO$ ), 3.9-3.65 (2H, m,  $NCHHCO$ ,  
 $NCHHCH_2$ ), 3.6-3.5 (1H, m,  $NCHHCH_2$ ), 2.83 (2H, t,  $^3J_{HH}$ =7Hz,  $NH_3CH_2$ ), 2.51 (2H, t,  
 $^3J_{HH}$ =7Hz,  $CH_2CO_2$ ), 2.17 (2H, t,  $^3J_{HH}$ =7Hz,  $C[CCH_2]$ ), 1.57 & 1.40 (each 2H, p,  
 $^3J_{HH}$ =7Hz,  $CH_2CH_2CH_2CH_2$ ).

10

## Example 8



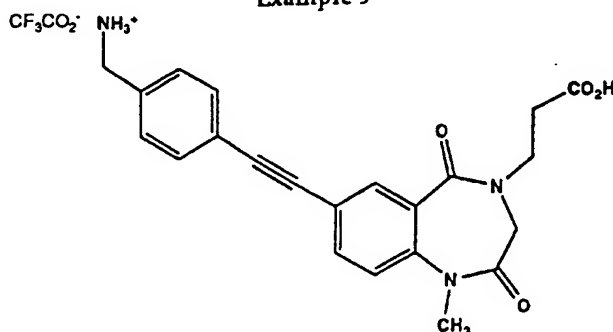
15 *1-(3-nitrobenzyl)-4-(2-carboxyethyl)-7-(6-guanidino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate.*

1-(3-nitrobenzyl)-4-(2-carboxyethyl)-7-(6-guanidino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate was prepared using the method described in Example 2. Thus, 45mgs of 1-(3-nitrobenzyl)-4-(2-carboxyethyl)-7-(6-amino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate yielded 25mgs (52%) of 1-(3-nitrobenzyl)-4-(2-carboxyethyl)-7-(6-guanidino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate (1/2" C-18 reverse-phase column, eluting with a solvent gradient of 20:80 methanol(0.1% trifluoroacetic acid)/ water (0.1% trifluoroacetic acid), time 0 to 10 min, to 70:30 methanol(0.1% trifluoroacetic acid)/water (0.1% trifluoroacetic acid), time 10 min to 40 min, flow=10ml/min,  $R_t$ =41.3 min, un detection 254nm). HRMS (FAB)  
 25 molecular ion  $m/z$ =521.2161 ( cald.  $C_{26}H_{28}N_6O_6$ , 521.2148).  $^1H$  NMR ( $D_2O$ ) 7.86 [1H, dt,  $^4J_{HH}$ =3Hz,  $^3J_{HH}$ =6Hz,  $NCH_2(C4-Ar-H)$ ], 7.64 [1H, bs,  $NCH_2(C2-Ar-H)$ ], 7.46 (1H, d,  $^4J_{HH}$ =2Hz,  $o$ -CON), 7.30-7.23 [3H, m,  $NCH_2(C5,C6-Ar-H)$ , Ar-H  $p$ -CON], 7.18 (1H, d,  $^3J_{HH}$ =8Hz, Ar-H  $m$ -CON), 5.30 (1H, d,  $^2J_{HH}$ =16Hz,  $NCHHAr$ ), 4.7 (1H,  $NCHHAr$ , overlapping with HOD), 4.03 (1H, d,  $^2J_{HH}$ =14Hz,  $NCHHCO$ ), 3.85 (1H, dt,  $^2J_{HH}$ =14Hz,

$^3J_{HH}=7\text{Hz}$ ,  $\text{NCHHCH}_2$ ), 3.74 (1H, d,  $^2J_{HH}=14\text{Hz}$ ,  $\text{NCHHCO}$ ), 3.55 (1H, dt,  $^2J_{HH}=14\text{Hz}$ ,  $^3J_{HH}=7\text{Hz}$ ,  $\text{NCHHCH}_2$ ), 2.97 (2H, t,  $^3J_{HH}=7\text{Hz}$ ,  $\text{H}_2\text{N}(\text{H}_2\text{N}=\text{N})\text{NHCH}_2$ ), 2.49 (2H, t,  $^3J_{HH}=7\text{Hz}$ ,  $\text{CH}_2\text{CO}_2$ ), 2.21 (2H, t,  $^3J_{HH}=7\text{Hz}$ ,  $\text{C}(\text{CCH}_2)$ ), 1.48 & 1.38 (each 2H, p,  $^3J_{HH}=7\text{Hz}$ ,  $\text{CH}_2\text{CH}_2\text{CH}_2\text{CH}_2$ ).

5

## Example 9



1-(methyl)-4-(2-carboxyethyl)-7-[4-(aminomethyl)phenyl]ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate.

- 10 a) 1-(methyl)-4-(2-carboxyethyl)-7-(4-cyanophenyl)ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester was prepared using the method employed in part (f) of Example 1. Thus, from 0.41 grams of 1-methyl-2-(carboxyethyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester and 0.127 grams of (4-cyanophenyl)ethyne (Gilbert, J.C.; Weerasooriya, U. *J.Org.Chem.* 1979, 44, 4997-98) was prepared 0.205 grams of 1-
- 15 (methyl)-4-(2-carboxyethyl)-7-(4-cyanophenyl)ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester (mp=160-161°C, TLC,  $\text{SiO}_2$ , 1:1 ethyl acetate/hexane,  $R_f=0.30$ , un positive).  $^1\text{H}$  NMR ( $\text{CDCl}_3$ , dTMS) 7.98 (1H, d,  $^4J_{HH}=2\text{Hz}$ , Ar-Ho-CON), 7.65-7.50 (5H, m, Ar-H p-CON, Ar-H o,p-CN), 7.17 (1H, d,  $^3J_{HH}=9\text{Hz}$ , m-CON), 4.09 (2H, q,  $^3J_{HH}=7\text{Hz}$ ,  $\text{OCH}_2$ ), 4.02 (1H, d,  $^2J_{HH}=15\text{Hz}$ ,  $\text{NCHHCO}$ ), 3.89 (2H, t,  $^3J_{HH}=7\text{Hz}$ ,  $\text{NCH}_2\text{CH}_2$ ), 3.86 (1H, d,  $^2J_{HH}=15\text{Hz}$ ,  $\text{NCHHCO}$ ), 3.34 (3H, s,  $\text{NCH}_3$ ), 2.72 (1H, dt,  $^2J_{HH}=16\text{Hz}$ ,  $^3J_{HH}=7\text{Hz}$ ,  $\text{CHHCO}_2$ ), 2.62 (1H, dt,  $^2J_{HH}=16\text{Hz}$ ,  $^3J_{HH}=7\text{Hz}$ ,  $\text{CHHCO}_2$ ), 1.20 (3H, t,  $^3J_{HH}=7\text{Hz}$ ,  $\text{CH}_3$ ).  $^{13}\text{C}$  NMR ( $\text{CDCl}_3$ , dTMS) 171.2, 168.6, 166.2, 141.0, 134.8, 134.4, 132.0, 128.8, 127.5, 121.1, 119.7, 118.3, 111.8, 91.7, 88.9, 60.3, 52.1, 45.1, 32.7, 31.5, 14.1.
- 20 b) A magnetically stirred solution of 3mL toluene, 6.8mgs dicobalt octacarbonyl (0.02 mmol) cooled to -20°C was saturated with triethylsilane *via* a gas bubbler. 0.1 grams of 1-(methyl)-4-(2-carboxyethyl)-7-(4-cyanophenyl)ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester (0.24 mmol) was added as a solution in 0.5mL toluene and the reaction mixture heated to 60°C for 20 hours (Murai, T. *et al.*, *Tetrahedron Lett.* 1985, 26, 5145-48). The
- 25 mixture was allowed to cool to room temperature and concentrated *in vacuo*. The resulting residue was diluted with methanol and 2mL of 2N sodium hydroxide was added. After 30 minutes, the mixture was neutralized with 2mL acetic acid, concentrated *in vacuo*, diluted

with 2mL of a 1:1 methanol/water solution and purified by high-pressure liquid chromatography (1/2" C-18 reverse-phase column, eluting with a solvent gradient of 20:80 methanol(0.1% trifluoroacetic acid)/ water (0.1% trifluoroacetic acid), time 0 to 10 min, to 70:30 methanol(0.1% trifluoroacetic acid)/water (0.1% trifluoroacetic acid), time 10 min to 40 min, flow=10ml/min,  $R_t$ =36.8 min, un detection 254nm) to yield 30.0 mgs (25%) of 1-(methyl)-4-(2-carboxyethyl)-7-[4-(aminomethyl)phenyl]ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate.  $^1\text{H}$  NMR ( $\text{CD}_3\text{OD}$ ) 7.68 (1H, d,  $^4\text{J}_{\text{HH}}=2\text{Hz}$ , Ar-H *o*-CON), 7.48 (1H, dd,  $^4\text{J}_{\text{HH}}=2\text{Hz}$ ,  $^3\text{J}_{\text{HH}}=9\text{Hz}$ , Ar-H *p*-CON) 7.4-7.3 (4H, m, Ar-H *o,p*-CN), 7.17 (1H, d, *m*-CON), 4.02 (2H, s,  $\text{ArCH}_2\text{N}$ ), 4.0-3.7 (4H, m,  $\text{NCH}_2\text{CO}$ ,  $\text{NCH}_2\text{CH}_2$ ), 3.0 (3H, s,  $\text{NCH}_3$ ) 2.6 (2H, m,  $\text{CH}_2\text{CO}_2$ ).

Using the above procedure, but substituting the appropriate 3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione nitrile for 1-(methyl)-4-(2-carboxyethyl)-7-(4-cyanophenyl)ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester there may be prepared, for example, the following compounds:

- 15 1-methyl-4-(2-carboxy-1-methylethyl)-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,  
1-methyl-4-(2-carboxy-1-phenylethyl)-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
( $\pm$ )-1-methyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 ( $\pm$ )-1-methyl-4-[succin-2-yl-(1-glycine)]-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
( $\pm$ )-1-methyl-4-[succin-2-yl-(1-valine)]-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 ( $\pm$ )-1-methyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-phenyl-4-(2-carboxy-1-methylethyl)-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-phenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 ( $\pm$ )-1-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
( $\pm$ )-1-phenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 ( $\pm$ )-1-phenyl-4-[succin-2-yl-(1-valine)]-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
( $\pm$ )-1-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- 1-benzyl-4-(2-carboxy-1-methylethyl)-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-benzyl-4-(2-carboxy-1-phenylethyl)-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1-benzyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-glycine)]-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-valine)]-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-
- 10 dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-isopropyl-4-(2-carboxy-1-methylethyl)-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 1-isopropyl-4-(2-carboxy-1-phenylethyl)-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-glycine)]-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-
- 20 dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-valine)]-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 1-methyl-4-(2-carboxy-1-methylethyl)-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-methyl-4-(2-carboxy-1-phenylethyl)-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-
- 30 dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-glycine)]-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-valine)]-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-
- 35 dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-phenyl-4-(2-carboxy-1-methylethyl)-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- 1-phenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1-phenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-valine)]-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 1-benzyl-4-(2-carboxy-1-methylethyl)-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-benzyl-4-(2-carboxy-1-phenylethyl)-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-glycine)]-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1-benzyl-4-[succin-2-yl-(1-valine)]-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-benzyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-isopropyl-4-(2-carboxy-1-methylethyl)-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 1-isopropyl-4-(2-carboxy-1-phenylethyl)-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-glycine)]-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 (±)-1-isopropyl-4-[succin-2-yl-(1-valine)]-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-isopropyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-methyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
(±)-1-methyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

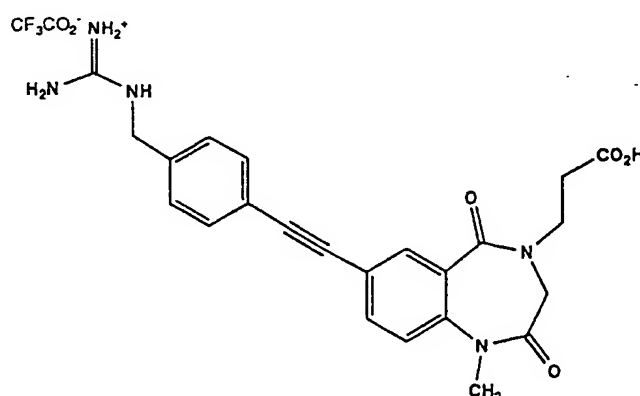
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-(2-carboxy-1-methylethyl)-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 (±)-1,3-diphenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1,3-diphenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-valine)]-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- 4-[succin-2-yl-(1-glycine)]-7-[2-(3-aminomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 (±)-1,3-diphenyl-4-(2-carboxy-1-methylethyl)-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1,3-diphenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-valine)]-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 (±)-1,3-diphenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(4-aminomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,



- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[2-(4-aminomethylthiophenol)]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(4-aminomethylthiophenol)]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 5 (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[2-(4-aminomethylthiophenol)]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(4-aminomethylthiophenol)]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(4-aminomethylthiophenol)]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 10 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(4-aminomethylthiophenol)]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[2-(4-aminomethylthiophenol)]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 15 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(4-aminomethylthiophenol)]-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate.

## Example 10



20

1-(methyl)-4-(2-carboxyethyl)-7-[4-(guanidinomethyl)phenyl]ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate.

- 1-(methyl)-4-(2-carboxyethyl)-7-[4-(guanidinomethyl)phenyl]ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate was prepared using the method described in  
 25 example 2. Thus, from 15 mgs of 1-(methyl)-4-(2-carboxyethyl)-7-[4-(aminomethyl)phenyl]ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate was prepared 7mgs (37%) of 1-(methyl)-4-(2-carboxyethyl)-7-[4-(guanidinomethyl)phenyl]ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate (1/2" C-18 reverse-phase column, eluting with a solvent gradient of 20:80  
 30 methanol(0.1% trifluoroacetic acid)/ water (0.1% trifluoroacetic acid), time 0 to 10 min, to

- 70:30 methanol(0.1% trifluoroacetic acid)/water (0.1% trifluoroacetic acid), time 10 min to 40 min, flow=10ml/min,  $R_t$ =39.6 min, un detection 254nm).  $^1\text{H}$  NMR ( $\text{CD}_3\text{OD}$ ,  $\text{D}_2\text{O}$ ) 7.50 (1H, d,  $^4\text{J}_{\text{HH}}=2\text{Hz}$ , Ar-H $\alpha$ -CON), 7.38 (1H, bd,  $^3\text{J}_{\text{HH}}=9\text{Hz}$ , Ar-H  $p$ -CON) 7.20 (1H, d,  $^3\text{J}_{\text{HH}}=9\text{Hz}$ ,  $m$ -CON), 7.15-7.0 (4H, m, Ar-H  $o,p$ -CN), 4.23 (2H, s,  $\text{ArCH}_2\text{NHC}(=\text{NH}_2)\text{NH}_2$ ), 3.9-3.8 (2H, m,  $\text{NCHHCO}$ ,  $\text{NCHHCH}_2$ ), 3.72 (1H, d,  $2\text{J}_{\text{HH}}=16\text{Hz}$ ,  $\text{NCHHCO}$ ), 3.58 (1H, m,  $\text{NCHHCH}_2$ ), 3.12 (3H, s,  $\text{NCH}_3$ ) 2.53 (2H, m,  $\text{CH}_2\text{CO}_2$ ).

- Using the above procedure, but substituting the appropriate amino acid for 1-(methyl)-4-(2-carboxyethyl)-7-[4-(aminomethyl)phenyl]ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate there may be prepared, for example, the following compounds:
- 1-methyl-4-(2-carboxy-1-methylethyl)-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione,
- 1-methyl-4-(2-carboxy-1-phenylethyl)-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- ( $\pm$ )-1-methyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- ( $\pm$ )-1-methyl-4-[succin-2-yl-(1-glycine)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- ( $\pm$ )-1-methyl-4-[succin-2-yl-(1-valine)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- ( $\pm$ )-1-methyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-phenyl-4-(2-carboxy-1-methylethyl)-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-phenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- ( $\pm$ )-1-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- ( $\pm$ )-1-phenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- ( $\pm$ )-1-phenyl-4-[succin-2-yl-(1-valine)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- ( $\pm$ )-1-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-benzyl-4-(2-carboxy-1-methylethyl)-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-benzyl-4-(2-carboxy-1-phenylethyl)-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1-benzyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-4-[succin-2-yl-(1-glycine)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1-benzyl-4-[succin-2-yl-(1-valine)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 1-isopropyl-4-(2-carboxy-1-methylethyl)-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-isopropyl-4-(2-carboxy-1-phenylethyl)-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1-isopropyl-4-[succin-2-yl-(1-glycine)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-4-[succin-2-yl-(1-valine)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 1-methyl-4-(2-carboxy-1-methylethyl)-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-methyl-4-(2-carboxy-1-phenylethyl)-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1-methyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-4-[succin-2-yl-(1-glycine)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-4-[succin-2-yl-(1-valine)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 (±)-1-methyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-phenyl-4-(2-carboxy-1-methylethyl)-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 1-phenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

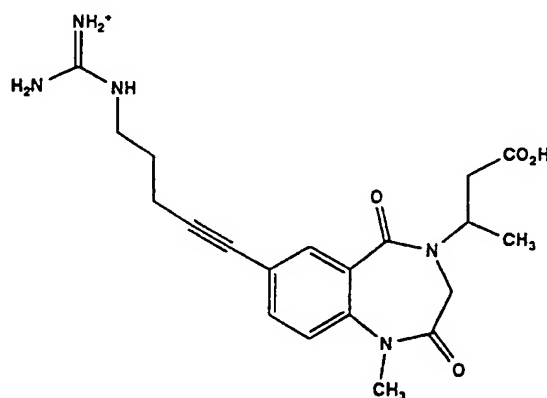
- (±)-1-phenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-phenyl-4-[succin-2-yl-(1-valine)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-benzyl-4-(2-carboxy-1-methylethyl)-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-benzyl-4-(2-carboxy-1-phenylethyl)-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 (±)-1-benzyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-4-[succin-2-yl-(1-glycine)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1-benzyl-4-[succin-2-yl-(1-valine)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-isopropyl-4-(2-carboxy-1-methylethyl)-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 1-isopropyl-4-(2-carboxy-1-phenylethyl)-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1-isopropyl-4-[succin-2-yl-(1-glycine)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-4-[succin-2-yl-(1-valine)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 (±)-1-methyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1,3-diphenyl-4-(2-carboxy-1-methylethyl)-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 (±)-1,3-diphenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-valine)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1,3-diphenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(3-guanidinomethylphenyl)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 (±)-1-methyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-methyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 (±)-1,3-diphenyl-4-(2-carboxy-1-methylethyl)-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 (±)-1,3-diphenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 (±)-1,3-diphenyl-4-[succin-2-yl-(1-valine)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1,3-diphenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 (±)-1-benzyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- (±)-1-benzyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-methylethyl)-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-(2-carboxy-1-phenylethyl)-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-*t*-butoxy)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-glycine)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-valine)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- (±)-1-isopropyl-3-phenyl-4-[succin-2-yl-(1-phenylalanine)]-7-[2-(4-guanidinomethylthiophenol)-1-ethyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate.

## Example 11



*1-methyl-4-(2-carboxy-1-methylethyl)-7-(5-guanidino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate.*

- a) Methyl N-(2-amino-5-iodobenzoyl)-3-amino-3-methylpropanate was prepared using the method described in part (b) of Example 1. Thus, 1.72 grams of 5-iodoisatoic anhydride and 1.73 grams of methyl 3-aminobutanoate p-tosylate yielded 0.65 grams (30%) methyl N-(2-amino-5-iodobenzoyl)-3-aminobutanoate (TLC, SiO<sub>2</sub>, 1:1 EtOAc/hexane, R<sub>f</sub>=0.63, unpositive).

b) Methyl N-(2-methylamino-5-iodobenzoyl)-3-amino-3-methylpropanoate was prepared using the method described in part (c) of Example 1. Thus, 0.32 grams of methyl N-(2-amino-5-iodobenzoyl)-3-aminobutanoate yielded 0.14 grams (43%) of methyl N-(2-methylamino-5-iodobenzoyl)-3-aminobutanoate (TLC, SiO<sub>2</sub>, 1:1 EtOAc/hexane, R<sub>f</sub>=0.85, un positive).

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c) 1-methyl-4-(3-methoxy-3-oxobutan-2-yl)-7-iodo-3,4-dihydro-1H-1,4-benzo-diazapine-2,5-dione was prepared using the method described in part (d) of Example 1. Thus, 0.25 grams of N-(2-methylamino-5-iodobenzoyl)-3-aminobutanoate yielded 0.08 grams (29%) of 1-methyl-4-(3-methoxy-3-oxobutan-2-yl)-7-iodo-3,4-dihydro-1H-1,4-benzo-diazapine-2,5-dione (TLC, SiO<sub>2</sub>, 1:1 EtOAc/hexane, R<sub>f</sub>=0.33, un positive).

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d) 1-methyl-4-(3-methoxy-3-oxobutan-2-yl)-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione was prepared using the method described in part (f) of Example 1. Thus, 0.08 grams of 1-methyl-4-(3-methoxy-3-oxobutan-2-yl)-7-iodo-3,4-dihydro-1H-1,4-benzo-diazapine-2,5-dione yielded 0.061 grams (68%) of 1-methyl-4-(3-methoxy-3-oxobutan-2-yl)-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione (TLC, SiO<sub>2</sub>, 1:1 EtOAc/hexane, R<sub>f</sub>=0.24, un positive).

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e) 1-methyl-4-(2-carboxy-1-methylethyl)-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate was prepared using the method described in part (g) of Example 1. Thus, 61mgs of 1-methyl-4-(3-methoxy-3-oxobutan-2-yl)-7-[5-(N-Boc)-amino-1-pentyne]-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione yielded 40 mgs (87%) of 1-methyl-4-(2-carboxy-1-methylethyl)-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate (1/2" C-18 reverse-phase column, eluting with a solvent gradient of 10:90 methanol(0.1% trifluoroacetic acid)/water(0.1% trifluoroacetic acid), time 0 to 10 min, to 50:50 methanol(0.1% trifluoroacetic acid)/water(0.1% trifluoroacetic acid), time 10 min to 40 min, flow 10ml/min, R<sub>t</sub>=31.2 min, un detection 254nm). HRMS (FAB) molecular ion m/z=358.1765 (cald. C<sub>19</sub>H<sub>24</sub>N<sub>3</sub>O<sub>4</sub>, 358.1767).

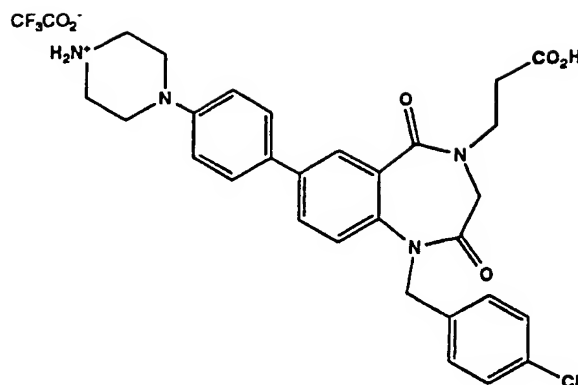
25

f) 1-methyl-4-(2-carboxy-1-methylethyl)-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate was prepared using the method described in Example 2. Thus, 15 mgs of 1-methyl-4-(2-carboxy-1-methylethyl)-7-(5-amino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate yielded 10mgs (67%) of 1-methyl-4-(2-carboxy-1-methylethyl)-7-(5-guanidino-1-pentyne)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate (1/2" C-18 reverse-phase column, eluting with a solvent gradient of 10:90 methanol(0.1% trifluoroacetic acid)/water(0.1% trifluoroacetic acid), time 0 to 10 min, to 50:50 methanol(0.1% trifluoroacetic acid)/water(0.1% trifluoroacetic acid), time 10 min to 40 min, flow 10ml/min, R<sub>t</sub>=36.9 min, un detection 254nm). HRMS (FAB) molecular ion m/z=400.1999 (cald. C<sub>20</sub>H<sub>26</sub>N<sub>5</sub>O<sub>4</sub>, 400.1986).

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## Example 12



5 **1-(4-chlorophenyl)methyl-4-(2-carboxyethyl)-7-(4-(1-piperidin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate.**

a) Over a period of 15 minutes samples of iodine (8.62 grams, 0.034mol) and silver benzoate (7.78 grams, 0.034mol) were added in portions to a magnetically stirred solution of 5 grams of 1-phenylpiperazine in 80mL acetic acid. The mixture stirred for 1 hour. The solvent was  
 10 evaporated and the residue partitioned between 150mL diethyl ether and 150mL 2N sodium hydroxide. The organic layer was washed with 100mL water, 100mL brine, then dried over potassium carbonate and concentrated *in vacuo* to yield 1.0 gram (11%) of 1-(4-iodophenyl)-piperzine (110-115°C). The material obtained was used without further purification. <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 7.51 (2H, d, <sup>3</sup>J<sub>HH</sub>=9Hz, *m*-ArH), 6.69 (2H, d, <sup>3</sup>J<sub>HH</sub>=9Hz, *o*-ArH), 3.17  
 15 (4H, m), 3.04 (4H, m). <sup>13</sup>C NMR (CDCl<sub>3</sub>) 151.36, 137.73, 118.05, 81.35, 49.92, 45.95.

b) The crude material obtained from part (a) (1 gram, 3.47mmol) was dissolved in dry THF (10mL) and a catalytic amount of dimethylamino-pyridine (100mgs) and di-*tert*-butyl dicarbonate (0.9 grams, 4.1mmol, 1.2 molar excess) were added, monitoring the reaction by  
 20 TLC (SiO<sub>2</sub>, 1:1 EtOAc/hexane, R<sub>f</sub>(product)=0.88). After 12 hours (overnight) the reaction was diluted with diethyl ether and extracted 1X75mL water, 1X75mL 1N HCl, 1X75mL sat. sodium bicarbonate, 1X75mL brine, dried over potassium carbonate, filtered and concentrated *in vacuo*. The material was further purified by filtering through a packet of SiO<sub>2</sub> using 1:1 diethyl ether/hexane as the eluting solvent to yield 1-(4-iodophenyl)-4-(N-Boc)-piperzine  
 25 (1 grm, 77%, 142-143°C). <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 7.54 (2H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, *m*-ArH), 6.69 (2H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, *o*-ArH), 3.58 (4H, t, <sup>3</sup>J<sub>HH</sub>=5Hz), 3.04 (4H, t, <sup>3</sup>J<sub>HH</sub>=5Hz) 1.48 (9H, s, Bu<sup>t</sup>).

c) A magnetically stirred solution of N-(2-amino-5-iodobenzoyl)-3-aminopropionate (8.8 grams, 23.4mmol), prepared using the procedure shown in part (b) of Example 1,  
 30 dimethylformamide, 2,6-lutidine (1.5 molar equiv., 35.1mmol, 4.25mL) and 4-chlorobenzyl chloride (1.5 molar equiv., 35.1mmol, 5.96 grams) was heated to 100°C for 24 hours. The

mixture was allowed to cool to room temperature and concentrated *in vacuo*. The resulting residue was dissolved in 150mL methylene chloride and washed with 2X100mL 10% citric acid, 1X100mL water, dried over sodium sulfate, filtered and concentrated *in vacuo*. The solid obtained was dissolved in a minimum volume of methylene chloride and a 5 fold excess of hexanes added. The solution was allowed to stand overnight in the refrigerator (5-7°C) to yield 7.56 grams of N-(2-(*p*-chlorobenzyl)amino-5-iodobenzoyl)-3-aminopropionate (153-155°C). Concentration of the mother liquor and repeating the steps for crystallization yielded a second crop (3.32 grams) for a total 92% isolated yield of the desired product. <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 8.10 (1H, bt, <sup>3</sup>J<sub>HH</sub>=5Hz, NH), 7.60 (1H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, C6-H), 7.44 (1H, dd, <sup>3</sup>J<sub>HH</sub>=9Hz, <sup>4</sup>J<sub>HH</sub>=2Hz, C8-H), 7.29 (4H, m, *p*-ClC<sub>6</sub>H<sub>4</sub>), 6.72 (1H, bt, <sup>3</sup>J<sub>HH</sub>=6Hz, NH), 6.34 (1H, d, <sup>3</sup>J<sub>HH</sub>=9Hz, C9-H), 4.35 (2H, d, <sup>3</sup>J<sub>HH</sub>=9Hz, NCH<sub>2</sub>Ar), 4.20 (2H, q, <sup>3</sup>J<sub>HH</sub>=7Hz, OCH<sub>2</sub>), 3.68 (2H, q, <sup>3</sup>J<sub>HH</sub>=6Hz, NCH<sub>2</sub>), 2.65 (2H, t, <sup>3</sup>J<sub>HH</sub>=6, CH<sub>2</sub>CO<sub>2</sub>), 1.31 (3H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, CH<sub>2</sub>CH<sub>3</sub>). <sup>13</sup>C NMR (CDCl<sub>3</sub>) 172.6, 168.3, 148.6, 141.0, 137.0, 135.6, 128.8, 128.3, 117.6, 114.3, 75.1, 60.9, 46.3, 35.1, 33.9, 14.2

d) To a magnetically stirred biphasic solution of N-(2-(*p*-chlorobenzyl)amino-5-iodobenzoyl)-3-aminopropionate (10grams, 20.57mmol), 200mL methylene chloride, and 200mL water was added, at room temperature, *o*-bromoacetyl bromide (24.7mmol, 1.2 molar equiv., 2.15mL), monitoring the reaction by TLC (SiO<sub>2</sub>, 1:1 EtOAc/hexane). The reaction was complete 6 hours. The layers were separated, washing the aqueous layer with 75mL methylene chloride. The combined organics were washed 1X100mL sat. sodium bicarbonate, 1X100mL Brine, dried over sodium sulfate, filtered and concentrated *in vacuo*. The resulting residue was dissolved in 50mL dimethylformamide and added, *via* an addition funnel, over a period of 30 minutes to a slurry of sodium hydride (1.2 molar equiv., 24.7mmol, 0.60 grams) in 10mL dimethylformamide cooled to 0°C under an atmosphere of nitrogen. The reaction was stirred for an additional 1 hour and then poured over an ice-cold 10% citric acid solution. The cloudy mixture was extracted 3X100mL EtOAc. The combined organics were dried over magnesium sulfate, filtered, and concentrated *in vacuo*. Purification was achieved by column chromatography (SiO<sub>2</sub>, 600mL, using 30%EtOAc/70% hexane to 60%EtOAc/40% hexane as the eluting solvent gradient) to yield 6.5 grams (63%) of 1-(*p*-chlorobenzyl)-7-iodo-4-(2-carboxyethyl)-3,4-dihydro-1*H*-1,4-benzodiazepine-2,4-dione ethyl ester. <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 8.11 (1H, d, <sup>3</sup>J<sub>HH</sub>=2Hz, C6-H), 6.68 (1H, dd, <sup>4</sup>J<sub>HH</sub>=2Hz, <sup>3</sup>J<sub>HH</sub>=9Hz, C8-H), 7.23 (2H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, ArH), 7.03 (2H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, ArH), 6.87 (1H, d, <sup>3</sup>J<sub>HH</sub>=9Hz, C9-H), 5.05 (1H, d, <sup>2</sup>J<sub>HH</sub>=16Hz, NCHHAr), 4.95 (1H, d, <sup>2</sup>J<sub>HH</sub>=16Hz, NCHHAr), 4.1 (3H, m, OCH<sub>2</sub>, C3-H), 3.93 (1H, d, <sup>2</sup>J<sub>HH</sub>=15Hz, C3-H), 3.9 (2H, m, NCH<sub>2</sub>CH<sub>2</sub>), 2.65 (2H, m, CH<sub>2</sub>CO<sub>2</sub>), 1.25 (3H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, OCH<sub>2</sub>CH<sub>3</sub>)

e) To a magnetically stirred solution of 1-(4-iodophenyl)-4-(*N*-Boc)-piperazine (0.5 grams, 1.29mmol) in 0.65mL THF at -78°C under an atmosphere of nitrogen was added *n*-BuLi as a

2.5M solution in hexane (1.42mmol, 1.1 molar equiv., 0.57mL) over a 5 minute period. The mixture was stirred for 15 minutes at -78°C and the treated with triisopropylborate (3 molar equiv., 3.9mmol, 0.9mL). The mixture was allowed to warm to room temperature and stirred overnight (ca. 12 hours). The mixture was cooled to 0°C and acidified to pH 6.5 with 5% aq. HCl and extracted with CH<sub>2</sub>Cl<sub>2</sub> (50mL). The organic layer was washed with brine, dried over sodium sulfate, decanted and concentrated *in vacuo*. The resulting residue, in a minimum volume of ethanol, was added to a prestirred (10 mins) mixture of tetrakis(triphenyl)phosphine palladium (0.03 molar equiv., 0.026mmol, 30mgs), and 0.447 grams of 1-(p-chlorobenzyl)-7-iodo-4-(2-carboxyethyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,4-dione ethyl ester (0.85mmol) in 0.5mL dimethylformamide at room temperature. 1.3mL of a 2M solution of sodium carbonate (2 equiv.) was added and the reaction mixture heated to 80°C for 6 hours. The mixture was allowed to cool to room temperature and is quenched with 1N HCl (25mL) and extracted 3X50mL ethyl acetate. The combined organics were washed 1X50mL sat NaHCO<sub>3</sub>, brine, dried over Na<sub>2</sub>SO<sub>4</sub>, decanted and concentrated *in vacuo*. The resulting residue was chromatographed (SiO<sub>2</sub>, 1:1 EtOAc/hexane to 9:1 EtOAc/hexane as the eluting solvent) to yield 1-(p-chlorobenzyl)-7-[4-(4-N-Boc-1-piperazine)phenyl]-4-(2-carboxyethyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,4-dione ethyl ester (0.137 grams, 24%, R<sub>f</sub> (SiO<sub>2</sub>, 1:1 EtOAc/hexane)=0.15, mu and ninhydrin positive). <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 8.00 (1H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, C6-H), 7.59 (1H, dd, <sup>4</sup>J<sub>HH</sub>=2Hz, <sup>3</sup>J<sub>HH</sub>=9Hz, C8-H), 7.50 (2H, d, <sup>3</sup>J<sub>HH</sub>=9Hz, C7-o-ArH), 7.23 (2H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, CH<sub>2</sub>-m-ArH), 7.16 (1H, d, <sup>3</sup>J<sub>HH</sub>=9Hz, C9-H), 7.10 (2H, d, <sup>3</sup>J<sub>HH</sub>=9Hz, CH<sub>2</sub>-o-ArH), 6.95 (2H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, C7-m-ArH), 5.08 (1H, d, <sup>2</sup>J<sub>HH</sub>=16Hz, NCHHAr), 4.99 (1H, d, <sup>2</sup>J<sub>HH</sub>=16Hz, NCHHAr), 4.18 (1H, d, <sup>2</sup>J<sub>HH</sub>=15Hz, C3-H), 4.13 (2H, q, <sup>3</sup>J<sub>HH</sub>=7Hz, OCH<sub>2</sub>), 3.92 (2H, m, NCH<sub>2</sub>CH<sub>2</sub>), 3.92 (1H, d, <sup>2</sup>J<sub>HH</sub>=15Hz, C3-H), 3.58 (4H, t, <sup>3</sup>J<sub>HH</sub>=5Hz, NCH<sub>2</sub>CH<sub>2</sub>NH), 3.18 (4H, t, <sup>3</sup>J<sub>HH</sub>=5Hz, NCH<sub>2</sub>NCH<sub>2</sub>NH), 2.68 (2H, m, CH<sub>2</sub>CO<sub>2</sub>), 1.47 (9H, s, Bu<sup>t</sup>), 1.25 (3H, t, <sup>3</sup>J<sub>HH</sub>=7Hz). <sup>13</sup>C NMR 171.3, 168.5, 167.1, 154.7, 150.9, 138.6, 138.0, 135.2, 133.3, 132.1, 130.9, 129.9, 129.7, 128.9, 128.4, 127.5, 121.9, 116.4, 79.9, 60.8, 52.4, 50.6, 48.8, 45.1, 32.8, 28.4, 14.2.

Using the above procedure, but substituting the appropriate 7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,4-dione alkyl ester for 1-(p-chlorobenzyl)-7-iodo-4-(2-carboxyethyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,4-dione ethyl ester and substituting the appropriate iodoarene for 1-(4-iodophenyl)-4-(N-Boc)-piperazine there may be prepared, for example, the following compounds:

- 1-methyl-7-[4-(4-N-Boc-1-piperazine)phenyl]-4-(2-carboxyethyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,4-dione ethyl ester,
- 1-(diphenylmethyl)-7-[4-(4-N-Boc-1-piperazine)phenyl]-4-(2-carboxyethyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,4-dione ethyl ester,
- 1-(2-naphthyl)methyl-7-[4-(4-N-Boc-1-piperazine)phenyl]-4-(2-carboxyethyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,4-dione ethyl ester,

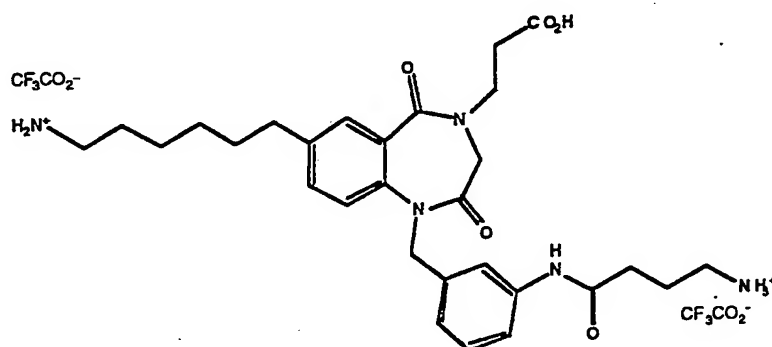
- (±)-1-methyl-7-[4-(4-N-Boc-1-piperazine)phenyl]-4-(3-phenyl-3-propionate)-3,4-dihydro-1H-1,4-benzodiazpine-2,4-dione methyl ester,  
 (±)-1-(2-naphthyl)methyl-7-[4-(4-N-Boc-1-piperazine)phenyl]-4-(3-phenyl-3-propionate)-3,4-dihydro-1H-1,4-benzodiazpine-2,4-dione methyl ester,
- 5 1-(p-chlorobenzyl)-7-[4-(2-(N-Boc)-aminoethoxy)phenyl]-4-(2-carboxyethyl)-3,4-dihydro-1H-1,4-benzodiazpine-2,4-dione ethyl ester  
 1-methyl-7-[4-(2-(N-Boc)-aminoethoxy)phenyl]-4-(2-carboxyethyl)-3,4-dihydro-1H-1,4-benzodiazpine-2,4-dione ethyl ester,  
 1-(diphenylmethyl)-7-[4-(2-(N-Boc)-aminoethoxy)phenyl]-4-(2-carboxyethyl)-3,4-dihydro-1H-1,4-benzodiazpine-2,4-dione ethyl ester,
- 10 1-(2-naphthyl)methyl-7-[4-(2-(N-Boc)-aminoethoxy)phenyl]-4-(2-carboxyethyl)-3,4-dihydro-1H-1,4-benzodiazpine-2,4-dione ethyl ester,  
 (±)-1-methyl-7-[4-(2-(N-Boc)-aminoethoxy)phenyl]-4-(3-phenyl-3-propionate)-3,4-dihydro-1H-1,4-benzodiazpine-2,4-dione methyl ester,
- 15 (±)-1-(2-naphthyl)methyl-7-[4-(2-(N-Boc)-aminoethoxy)phenyl]-4-(3-phenyl-3-propionate)-3,4-dihydro-1H-1,4-benzodiazpine-2,4-dione methyl ester,  
 1-(p-chlorobenzyl)-7-[4-(N-Boc-aminomethyl)phenyl]-4-(2-carboxyethyl)-3,4-dihydro-1H-1,4-benzodiazpine-2,4-dione ethyl ester  
 1-methyl-7-[4-(N-Boc-aminomethyl)phenyl]-4-(2-carboxyethyl)-3,4-dihydro-1H-1,4-benzodiazpine-2,4-dione ethyl ester,
- 20 1-(diphenylmethyl)-7-[4-(N-Boc-aminomethyl)phenyl]-4-(2-carboxyethyl)-3,4-dihydro-1H-1,4-benzodiazpine-2,4-dione ethyl ester,  
 1-(2-naphthyl)methyl-7-[4-(N-Boc-aminomethyl)phenyl]-4-(2-carboxyethyl)-3,4-dihydro-1H-1,4-benzodiazpine-2,4-dione ethyl ester,
- 25 (±)-1-methyl-7-[4-(N-Boc-aminomethyl)phenyl]-4-(3-phenyl-3-propionate)-3,4-dihydro-1H-1,4-benzodiazpine-2,4-dione methyl ester,  
 (±)-1-(2-naphthyl)methyl-7-[4-(N-Boc-aminomethyl)phenyl]-4-(3-phenyl-3-propionate)-3,4-dihydro-1H-1,4-benzodiazpine-2,4-dione methyl ester.
- 30 f) 1-(p-chlorobenzyl)-7-[4-(1-piperazine)phenyl]-4-(2-carboxyethyl)-3,4-dihydro-1H-1,4-benzodiazpine-2,4-dione trifluoroacetate was prepared using the method described in part (c) of Example 1. Thus, 34 mgs (0.052mmol) 1-(p-chlorobenzyl)-7-[4-(4-N-Boc-1-piperazine)phenyl]-4-(2-carboxyethyl)-3,4-dihydro-1H-1,4-benzodiazpine-2,4-dione ethyl ester yielded 25 mgs (90%) of 1-(p-chlorobenzyl)-7-[4-(1-piperazine)phenyl]-4-(2-carboxyethyl)-3,4-dihydro-1H-1,4-benzodiazpine-2,4-dione trifluoroacetate (1/2" C-18 reverse-phase column, eluting with a solvent gradient of 1:9 acetonitrile(0.1%TFA)/water (0.1% TFA) time 0 to 10 minutes, to 1:1 acetonitrile(0.1%TFA)/water (0.1% TFA) time 10 to 40 minutes, flow 10mL/min,  $R_t$ =34.9 min, mu detection 254nm). HRMS (FAB) molecular ion  $m/z$ =533.1950 (calcd.  $C_{29}H_{30}N_4O_4Cl$ , 533.1956).  $^1H$  NMR ( $D_2O$ , 10% $CD_3OD$ ) 7.61 (1H, d,
- 35

- <sup>4</sup>J<sub>HH</sub>=2Hz, C6-H), 7.43 (1H, dd, <sup>4</sup>J<sub>HH</sub>=2Hz, <sup>3</sup>J<sub>HH</sub>=8Hz, C8-H), 7.28 (2H, d, C7-*o*-ArH), 7.10 (1H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, C9-H), 6.98 (2H, d, <sup>3</sup>J<sub>HH</sub>=7Hz, CH<sub>2</sub>-*m*-ArH), 6.87 (2H, d, <sup>3</sup>J<sub>HH</sub>=7Hz, CH<sub>2</sub>-*o*-ArH), 6.80 (2H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, C7-*m*-ArH), 5.20 (1H, d, <sup>2</sup>J<sub>HH</sub>=16Hz, NCHHAr), 3.95 (1H, d, <sup>2</sup>J<sub>HH</sub>=14Hz, C3-H), 3.84 (1H, dt, NCHHCH<sub>2</sub>), 3.72 (1H, d, <sup>2</sup>J<sub>HH</sub>=14Hz, C3-H), 3.58 (1H, dt, C3-H), 3.23 (4H, bs, NCH<sub>2</sub>CH<sub>2</sub>NH), 3.18 (4H, bs, NCH<sub>2</sub>CH<sub>2</sub>NH), 2.58 (2H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, CH<sub>2</sub>CO<sub>2</sub>).

- Using the above procedure, but substituting the appropriate 7-aryl-3,4-dihydro-1*H*-1,4-benzodiazepine-2,4-dione N-Boc amino alkyl ester for 1-(*p*-chlorobenzyl)-7-[4-(4-N-Boc-1-piperazine)phenyl]-4-(2-carboxyethyl)-3,4-dihydro-1*H*-1,4-benzodiazepine-2,4-dione ethyl ester there may be prepared, for example, the following compounds:
- 1-methyl-7-[4-(1-piperazine)phenyl]-4-(2-carboxyethyl)-3,4-dihydro-1*H*-1,4-benzodiazepine-2,4-dione trifluoroacetate,
- 1-(diphenylmethyl)-7-[4-(1-piperazine)phenyl]-4-(2-carboxyethyl)-3,4-dihydro-1*H*-1,4-benzodiazepine-2,4-dione trifluoroacetate,
- 1-(2-naphthyl)methyl-7-[4-(1-piperazine)phenyl]-4-(2-carboxyethyl)-3,4-dihydro-1*H*-1,4-benzodiazepine-2,4-dione trifluoroacetate,
- (±)-1-methyl-7-[4-(1-piperazine)phenyl]-4-(3-phenyl-3-proprionate)-3,4-dihydro-1*H*-1,4-benzodiazepine-2,4-dione trifluoroacetate,
- (±)-1-(2-naphthyl)methyl-7-[4-(1-piperazine)phenyl]-4-(3-phenyl-3-proprionate)-3,4-dihydro-1*H*-1,4-benzodiazepine-2,4-dione trifluoroacetate,
- 1-(*p*-chlorobenzyl)-7-[4-(2-aminoethoxy)phenyl]-4-(2-carboxyethyl)-3,4-dihydro-1*H*-1,4-benzodiazepine-2,4-dione trifluoroacetate,
- 1-methyl-7-[4-(2-aminoethoxy)phenyl]-4-(2-carboxyethyl)-3,4-dihydro-1*H*-1,4-benzodiazepine-2,4-dione trifluoroacetate,
- 1-(diphenylmethyl)-7-[4-(2-aminoethoxy)phenyl]-4-(2-carboxyethyl)-3,4-dihydro-1*H*-1,4-benzodiazepine-2,4-dione trifluoroacetate,
- 1-(2-naphthyl)methyl-7-[4-(2-aminoethoxy)phenyl]-4-(2-carboxyethyl)-3,4-dihydro-1*H*-1,4-benzodiazepine-2,4-dione trifluoroacetate,
- (±)-1-methyl-7-[4-(2-aminoethoxy)phenyl]-4-(3-phenyl-3-proprionate)-3,4-dihydro-1*H*-1,4-benzodiazepine-2,4-dione trifluoroacetate,
- (±)-1-(2-naphthyl)methyl-7-[4-(2-aminoethoxy)phenyl]-4-(3-phenyl-3-proprionate)-3,4-dihydro-1*H*-1,4-benzodiazepine-2,4-dione trifluoroacetate,
- 1-(*p*-chlorobenzyl)-7-[4-(aminomethyl)phenyl]-4-(2-carboxyethyl)-3,4-dihydro-1*H*-1,4-benzodiazepine-2,4-dione trifluoroacetate,
- 1-methyl-7-[4-(aminomethyl)phenyl]-4-(2-carboxyethyl)-3,4-dihydro-1*H*-1,4-benzodiazepine-2,4-dione trifluoroacetate,
- 1-(diphenylmethyl)-7-[4-(aminomethyl)phenyl]-4-(2-carboxyethyl)-3,4-dihydro-1*H*-1,4-benzodiazepine-2,4-dione trifluoroacetate,

- 1-(2-naphthyl)methyl-7-[4-(aminomethyl)phenyl]-4-(2-carboxyethyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,4-dione trifluoroacetate,  
 (±)-1-methyl-7-[4-(aminomethyl)phenyl]-4-(3-phenyl-3-propionate)-3,4-dihydro-1H-1,4-benzodiazepine-2,4-dione trifluoroacetate,  
 5 (±)-1-(2-naphthyl)methyl-7-[4-(aminomethyl)phenyl]-4-(3-phenyl-3-propionate)-3,4-dihydro-1H-1,4-benzodiazepine-2,4-dione trifluoroacetate.

## Example 13



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1-[3-(N-(4-aminobutyryl)aminophenyl)methyl]-4-(2-carboxyethyl)-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione bistrifluoroacetate.

- a) 1-(3-nitrophenyl)methyl-4-(2-carboxyethyl)-7-(N-Boc-6-aminohexynyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione ethyl ester (110mgs, 0.18mmol) was reduced under an atmosphere of nitrogen in 4mL ethyl acetate in the presence of a catalytic amount of 10% Pd on carbon (25mgs), monitoring the reaction by TLC (SiO<sub>2</sub>, 1:1 ethyl acetate/hexane, R<sub>f</sub>(starting material)=0.33, R<sub>f</sub>(product)=0.15, mu and ninhydrin positive). After 3 hours, the reaction was filtered through Celite®, washing with ethyl acetate, and concentrated *in vacuo* to yield 97mgs (92%) of 1-(3-aminophenyl)methyl-4-(2-carboxyethyl)-7-(N-Boc-6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione ethyl ester. 1H NMR (CDCl<sub>3</sub>, dTMS) 7.58 (1H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, C6-H), 7.18 (1H, dd, <sup>4</sup>J<sub>HH</sub>=2Hz, <sup>3</sup>J<sub>HH</sub>=8Hz, C8-H), 7.09 (1H, d, <sup>3</sup>J<sub>HH</sub>=9Hz, CH<sub>2</sub>ArH), 7.04 (1H, t, <sup>3</sup>J<sub>HH</sub>=8Hz, CH<sub>2</sub>ArH), 6.54 (1H, bd, CH<sub>2</sub>ArH), 6.48 (1H, bs, CH<sub>2</sub>ArH), 4.99 (1H, d, <sup>2</sup>J<sub>HH</sub>=16Hz, CHHAr), 4.90 (1H, d, <sup>2</sup>J<sub>HH</sub>=16Hz, CHHAr), 4.53 (1H, vbs, NHBoc), 4.20-3.96 (4H, m, OCH<sub>2</sub>, C3-H, NCHHCH<sub>2</sub>), 3.92-3.80 (2H, m, C3-H, NCHHCH<sub>2</sub>), 3.55 (2H, bs, NH<sub>2</sub>), 3.06 (2H, bq, <sup>3</sup>J<sub>HH</sub>=6Hz, CH<sub>2</sub>NHBoc), 2.72 (2H, m, CH<sub>2</sub>CO<sub>2</sub>), 2.56 (2H, t, <sup>3</sup>J<sub>HH</sub>=7Hz), 1.66-1.20 (20H, m, (CH<sub>2</sub>)<sub>4</sub>, Bu<sup>t</sup>, OCH<sub>2</sub>CH<sub>3</sub>). 13C NMR (CDCl<sub>3</sub>) 171.4, 168.5, 167.4, 146.3, 140.7, 138.1, 137.9, 132.2, 130.1, 129.6, 129.0, 121.6, 117.4, 114.5, 113.6, 79.0, 60.8, 52.4, 51.1, 45.1, 40.5, 34.9, 32.8, 30.8, 29.9, 28.8, 28.4, 26.5, 14.2.

- 30 Using the above procedure, but substituting the appropriate 7-alkyl or aryl-3,4-dihydro-1H-1,4-benzodiazapine-2,4-dione N-Boc amino alkyl ester for 1-(3-nitrophenyl)methyl-4-(2-carboxyethyl)-7-(N-Boc-6-aminohexynyl)-3,4-dihydro-1H-1,4-

benzodiazapine-2,5-dione ethyl ester there may be prepared, for example, the following compounds:

- 1-(2-aminophenyl)methyl-4-(2-carboxyethyl)-7-(N-Boc-6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione ethyl ester,
- 5 1-(4-aminophenyl)methyl-4-(2-carboxyethyl)-7-(N-Boc-6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione ethyl ester,
- 1-(3-aminophenyl)methyl-4-(2-carboxyethyl)-7-[4-(4-N-Boc-1-piperazine)phenyl]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione ethyl ester,
- 1-(3-aminophenyl)methyl-4-(3-phenyl-3-proprionate)-7-[4-(4-N-Boc-1-
- 10 piperazine)phenyl]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione ethyl ester,
- 1-(3-aminophenyl)methyl-4-(3-phenyl-3-proprionate)-7-(N-Boc-6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione ethyl ester,

- b) A solution of the aniline 1-(3-aminophenyl)methyl-4-(2-carboxyethyl)-7-(N-Boc-6-
- 15 aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione ethyl ester (50mgs, 0.09mmol) and 0.5mL dimethylacetamide was added to a prestirred solution of N-Boc-4-aminobutyric acid (0.16mmol, 32.5mgs), 0.5mL dimethylacetamide, 0.5mL pyridine, and benzotriazol-1-yl-oxy-tris-(dimethylamino)phosphonium hexafluorophosphate (BOP Reagent) and a
- 20 catalytic amount of dimethylaminopyridine (DMAP, 5mgs), monitoring by TLC (8:2 ethyl acetate/hexane). After 6 hours, the reaction was poured over 50mL 1N HCl and extracted 2X50mL ethyl acetate. The combined organics was dried over sodium sulfate, decanted, and concentrated in vacuo. The resulting residue was purified by column chromatography, SiO<sub>2</sub>, using 7:3 ethyl acetate/hexane to 9:1 ethyl acetate/hexane, R<sub>f</sub>(8:2 ethyl acetate/hexane, product)=0.15, to yield 59mgs (86%) of 1-[3-(4-N-Boc-aminobutyl)aminophenyl]methyl-4-
- 25 (2-carboxyethyl)-7-(N-Boc-6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione ethyl ester. <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 8.78 (1H, bs, ArNHCO), 7.56 (1H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, C6-H), 7.44 (1H, bd, <sup>3</sup>J<sub>HH</sub>=8Hz, C8-H), 7.36 (1H, bs, CH<sub>2</sub>ArH), 7.17 (2H, m, CH<sub>2</sub>ArH), 7.09 (1H, t, <sup>3</sup>J<sub>HH</sub>=8Hz, CH<sub>2</sub>ArH), 6.81 (1H, bd, <sup>3</sup>J<sub>HH</sub>=8Hz, C9-H), 5.1-4.9 (3H, m, NHBoc, CH<sub>2</sub>Ar), 4.58 (1H, bs, NHBoc), 4.20-4.05 (4H, m, OCH<sub>2</sub>, C3-H, NCHHCH<sub>2</sub>), 3.87 (1H, d, <sup>2</sup>J<sub>HH</sub>=15Hz, C3-H), 3.80 (1H, dt, NCHHCH<sub>2</sub>), 3.16 (2H, bq, <sup>3</sup>J<sub>HH</sub>=6Hz, CH<sub>2</sub>NHBoc), 3.04 (2H, bq, <sup>3</sup>J<sub>HH</sub>=6Hz, CH<sub>2</sub>NHBoc), 2.71 (2H, m, CH<sub>2</sub>CO<sub>2</sub>), 2.54 (2H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, NCOCH<sub>2</sub>), 2.32 (2H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, ArCH<sub>2</sub>), 1.82 (2H, p, <sup>3</sup>J<sub>HH</sub>=7Hz, NCOCH<sub>2</sub>CH<sub>2</sub>), 1.6-1.2 (20H, m, (CH<sub>2</sub>)<sub>4</sub>, Bu<sup>t</sup>, OCH<sub>2</sub>CH<sub>3</sub>).
- 30

- Using the above procedure, but substituting the appropriate 7-alkyl or aryl-3,4-
- 35 dihydro-1H-1,4-benzodiazapine-2,4-dione N-Boc amino alkyl ester for 1-(3-aminophenyl)methyl-4-(2-carboxyethyl)-7-(N-Boc-6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione ethyl ester there may be prepared, for example, the following compounds:

- 1-[2-(4-N-Boc-aminobutyl)aminophenyl]methyl-4-(2-carboxyethyl)-7-(N-Boc-6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione ethyl ester,  
 1-(4-(4-N-Boc-aminobutyl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(N-Boc-6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione ethyl ester,  
 5 1-(3-(4-N-Boc-aminobutyl)aminophenyl)methyl-4-(2-carboxyethyl)-7-[4-(4-N-Boc-1-piperazine)phenyl]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione ethyl ester,  
 1-(3-(4-N-Boc-aminobutyl)aminophenyl)methyl-4-(3-phenyl-3-propionate)-7-[4-(4-N-Boc-1-piperazine)phenyl]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione ethyl ester,  
 1-(3-(4-N-Boc-aminobutyl)aminophenyl)methyl-4-(3-phenyl-3-propionate)-7-(N-Boc-6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione ethyl ester,  
 10 1-[3-(3-N-Boc-aminopropionyl)aminophenyl]methyl-4-(2-carboxyethyl)-7-(N-Boc-6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione ethyl ester,  
 1-[2-(3-N-Boc-aminopropionyl)aminophenyl]methyl-4-(2-carboxyethyl)-7-(N-Boc-6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione ethyl ester,  
 15 1-(4-(3-N-Boc-aminopropionyl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(N-Boc-6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione ethyl ester,  
 1-(3-(3-N-Boc-aminopropionyl)aminophenyl)methyl-4-(2-carboxyethyl)-7-[4-(4-N-Boc-1-piperazine)phenyl]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione ethyl ester,  
 1-(3-(3-N-Boc-aminopropionyl)aminophenyl)methyl-4-(3-phenyl-3-propionate)-7-[4-(4-N-Boc-1-piperazine)phenyl]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione ethyl ester,  
 20 1-(3-(3-N-Boc-aminopropionyl)aminophenyl)methyl-4-(3-phenyl-3-propionate)-7-(N-Boc-6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione ethyl ester.

- c) 1-[3-(N-(4-aminobutyl))aminophenyl]methyl-4-(2-carboxyethyl)-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione trifluoroacetate was prepared using the method described in part (c) of Example 1. Thus, 59 mgs (0.074mmol) 1-[3-(4-N-Boc-aminobutyl)aminophenyl]methyl-4-(2-carboxyethyl)-7-(N-Boc-6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione ethyl ester yielded 50 mgs (99%) of 1-[3-(N-(4-aminobutyl))aminophenyl]methyl-4-(2-carboxyethyl)-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione trifluoroacetate (1/2" C-18 reverse-phase column, eluting with a solvent gradient of 1:9 acetonitrile(0.1%TFA)/water (0.1% TFA) time 0 to 10 minutes, to 1:1 acetonitrile(0.1%TFA)/water (0.1% TFA) time 10 to 40 minutes, flow 10mL/min,  $R_t=30.5$  min, mu detection 254nm). HRMS (FAB) molecular ion  $m/z=538.3032$  (calcd.  $C_{29}H_{39}N_5O_5$ , 538.3030).  $^1H$  NMR ( $D_2O$ ) 7.21 (1H, bs, C6-H), 7.05-6.85 (5H, m,  $CH_2ArH$ , C8-H), 6.52 (1H, bd,  $^3J_{HH}=8Hz$ , C9-H), 5.03 (1H, d,  $^2J_{HH}=16Hz$ ,  $CHHAr$ ), 4.47 (1H, d,  $^2J_{HH}=16Hz$ ,  $CHHAr$ ), 3.86 (1H, d,  $^2J_{HH}=15Hz$ , C3-H), 3.80 (1H, dt,  $^3J_{HH}=7Hz$ ,  $^2J_{HH}=15Hz$ ,  $NCHHCH_2$ ), 3.63 (1H, d,  $^2J_{HH}=15Hz$ , C3-H), 3.51 (1H, dt,  $^3J_{HH}=7Hz$ ,  $^2J_{HH}=15Hz$ ,  $NCHHCH_2$ ), 2.83 (2H, t,  $^3J_{HH}=8Hz$ ,  $CH_2NH_2$ ), 2.68 (2H, t,  $^3J_{HH}=8Hz$ ,



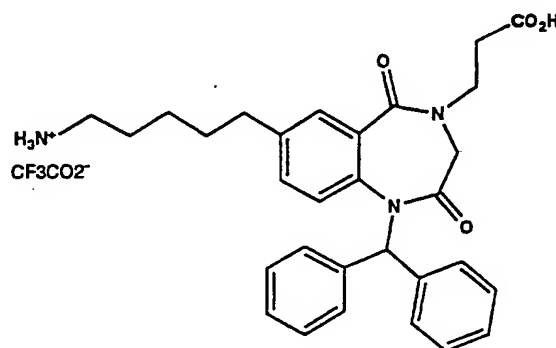
CH<sub>2</sub>NH<sub>2</sub>), 2.43 (2H, t, <sup>3</sup>J<sub>HH</sub>=7Hz), 2.3-2.2 (4H, m), 1.76 (2H, p, <sup>3</sup>J<sub>HH</sub>=7Hz, NCOCH<sub>2</sub>CH<sub>2</sub>), 1.35 (2H, p, <sup>3</sup>J<sub>HH</sub>=7Hz), 1.22 (2H, p, <sup>3</sup>J<sub>HH</sub>=7Hz), 1.1-0.9 (4H, m).

Using the above procedure, but substituting the appropriate 7-alkyl or aryl-3,4-dihydro-1H-1,4-benzodiazapine-2,4-dione bis-N-Boc diamino alkyl ester for 1-[3-(4-N-Boc-aminobutyryl)aminophenyl]methyl-4-(2-carboxyethyl)-7-(N-Boc-6-aminoheptyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione ethyl ester there may be prepared, for example, the following compounds:

- 1-[2-(4-aminobutyryl)aminophenyl]methyl-4-(2-carboxyethyl)-7-(6-aminoheptyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione bistrifluoroacetate,
- 10 1-(4-(4-aminobutyryl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(6-aminoheptyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione bistrifluoroacetate,
- 1-(3-(4-aminobutyryl)aminophenyl)methyl-4-(2-carboxyethyl)-7-[4-(1-piperazine)phenyl]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione bistrifluoroacetate,
- 1-(3-(4-aminobutyryl)aminophenyl)methyl-4-(3-phenyl-3-propionate)-7-[4-(1-piperazine)phenyl]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione bistrifluoroacetate,
- 15 1-(3-(4-aminobutyryl)aminophenyl)methyl-4-(3-phenyl-3-propionate)-7-(6-aminoheptyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione bistrifluoroacetate,
- 1-[3-(3-aminopropionyl)aminophenyl]methyl-4-(2-carboxyethyl)-7-(6-aminoheptyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione bistrifluoroacetate,
- 20 1-[2-(3-aminopropionyl)aminophenyl]methyl-4-(2-carboxyethyl)-7-(6-aminoheptyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione bistrifluoroacetate,
- 1-(4-(3-aminopropionyl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(6-aminoheptyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione bistrifluoroacetate,
- 1-(3-(3-aminopropionyl)aminophenyl)methyl-4-(2-carboxyethyl)-7-[4-(1-piperazine)phenyl]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione bistrifluoroacetate,
- 25 1-(3-(3-aminopropionyl)aminophenyl)methyl-4-(3-phenyl-3-propionate)-7-[4-(1-piperazine)phenyl]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione bistrifluoroacetate,
- 1-(3-(3-aminopropionyl)aminophenyl)methyl-4-(3-phenyl-3-propionate)-7-(6-aminoheptyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione bistrifluoroacetate.

30

#### Example 14



**1-(diphenylmethyl)-4-(2-carboxyethyl)-7-(5-aminopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate**

- 5 a) A magnetically stirred solution of 1.1 gram of N-(2-amino-5-iodobenzoyl)-b-alanine ethyl ester (3.1mmol), prepared by the method shown in part (b) of Example 1, 0.48mL of 2,6-lutidine (4.1mmol), 1.2 grams of chlorodiphenylmethane (4.7mmol), and 10mL of dimethylformamide was heated to 50°C for 1 hour. The reaction mixture was allowed to cool to room temperature and concentrated *in vacuo*. The resulting oil was dissolved in 35mL of
- 10 methylene chloride and washed with 2x50mL 10% citric acid, 1x50mL water, dried over sodium sulfate, decanted and concentrated *in vacuo*. The resulting oil was further purified by column chromatography, using silica gel, eluting with a solvent gradient of 10/90 ethyl acetate/hexane to 25/75 ethyl acetate/hexane (TLC, SiO<sub>2</sub>, 1:1 EtOAc/hexane, R<sub>f</sub> = 0.82, uv positive) to yield 0.92 gram (56%) of N-(2-diphenylmethylamino-5-iodobenzoyl)-b-alanine
- 15 ethyl ester. <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 8.40 (1H, d, <sup>3</sup>J<sub>HH</sub>=5Hz, NHCHAr<sub>2</sub>), 7.56 (1H, d, <sup>3</sup>J<sub>HH</sub>=2Hz, ArH o-CON), 7.18-7.36 (11H, m, <sup>4</sup>J<sub>HH</sub>=2Hz, ArH Ph, ArH p-CON), 6.73 (1H, t, <sup>3</sup>J<sub>HH</sub>=6Hz, NHCH<sub>2</sub>), 6.29 (1H, d, <sup>3</sup>J<sub>HH</sub>=9Hz, ArH m-CON), 5.52 (1H, d, <sup>3</sup>J<sub>HH</sub>=5Hz, CHNH), 4.15 (2H, q, <sup>3</sup>J<sub>HH</sub>=7Hz, OCH<sub>2</sub>), 3.61 (2H, q, <sup>3</sup>J<sub>HH</sub>=6Hz, CH<sub>2</sub>NH), 2.58 (2H, t, <sup>3</sup>J<sub>HH</sub>=6Hz, CH<sub>2</sub>CO), 1.25 (3H, t, <sup>3</sup>J<sub>HH</sub>=6Hz, CH<sub>3</sub>CH<sub>2</sub>O).

- 20 Using the above procedure, but substituting the appropriate alkyl or aryl halide there may be prepared, for example, the following compounds:
- N-(2-(1-naphthyl)methylamino-5-iodobenzoyl)-b-alanine ethyl ester,  
 N-(2-(2-naphthyl)methylamino-5-iodobenzoyl)-b-alanine ethyl ester,  
 N-(2-(p-phenoxybenzyl)amino-5-iodobenzoyl)-b-alanine ethyl ester,  
 25 N-(2-(m-phenoxybenzyl)amino-5-iodobenzoyl)-b-alanine ethyl ester,  
 N-(2-(p-trifluoromethylbenzyl)amino-5-iodobenzoyl)-b-alanine ethyl ester,  
 N-(2-(ethyl-5-valeroyl)amino-5-iodobenzoyl)-b-alanine ethyl ester,  
 N-(2-(p-methoxybenzyl)amino-5-iodobenzoyl)-b-alanine ethyl ester,  
 Methyl 3(S)-N-(2-(2-methylnaphthyl)amino-5-iodobenzoyl)-3-amino-3-methylpropionate,  
 30 Methyl 3(R)-N-(2-(2-naphthyl)methylamino-5-iodobenzoyl)-3-amino-3-methylpropionate,  
 Methyl (+/-)-N-(2-(2-naphthyl)methylamino-5-iodobenzoyl)-3-amino-3-methylpropionate,  
 Methyl 3(S)-N-(2-(2-naphthyl)methylamino-5-iodobenzoyl)-3-amino-3-phenylpropionate,

Methyl 3(R)-N-(2-(2-naphthyl)methylamino-5-iodobenzoyl)-3-amino-3-phenylpropionate,  
Methyl (+/-)-N-(2-(2-naphthyl)methylamino-5-iodobenzoyl)-3-amino-3-phenylpropionate,

- b) To a magnetically stirred solution of 0.93 gram of N-(2-diphenyl-methylamino-5-iodobenzoyl)-b-alanine ethyl ester (1.8mmol) in 15mL of methylene chloride and 15mL water, was added 183mL a-bromoacetyl bromide (2.1mmol). The reaction mixture was stirred overnight. The layers were separated, the aqueous fraction was washed 1x15mL methylene chloride, and the two methylene chloride fractions were combined and washed with 1x20mL of saturated sodium bicarbonate, dried over sodium sulfate, filtered and concentrated *in vacuo*. The resulting residue was dissolved in 7mL of dimethylformamide and added, *via* an addition funnel, to a slurry of 57 mgs of 95% sodium hydride (2.2mmol) in 5mL dimethylformamide that was cooled at 0°C. After 2 hours, the mixture was poured over 40mL of ice cooled 10% citric acid and extracted 3x40mL methylene chloride. The combined organic layers were washed with 2x50mL of 10% citric acid and 1x50mL water, dried over sodium sulfate, decanted, concentrated *in vacuo* and further purified by column chromatography, using silica gel, eluting with a 1:1 mixture of ethyl acetate and hexane (TLC, SiO<sub>2</sub>, 1:1 EtOAc/hexane, R<sub>f</sub> = 0.61, uv positive) to yield 474 mgs (48%) of 1-(diphenylmethyl)-4-(2-carboxyethyl)-8-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester. <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 8.03 (1H, d, <sup>3</sup>J<sub>HH</sub>=2Hz, ArH o-CON), 7.45 (1H, dd, <sup>3</sup>J<sub>HH</sub>=2Hz, <sup>3</sup>J<sub>HH</sub>=9Hz, ArH p-CON), 7.1-7.4 (1H, m, <sup>4</sup>J<sub>HH</sub>=2Hz, ArH Ph), 6.79 (1H, d, <sup>3</sup>J<sub>HH</sub>=9Hz, ArH m-CON), 6.70 (1H, s), 4.13 (2H, q, <sup>3</sup>J<sub>HH</sub>=7Hz, OCH<sub>2</sub>), 3.98 (2H, dd, <sup>3</sup>J<sub>HH</sub>=100Hz, <sup>3</sup>J<sub>HH</sub>=15Hz, COCH<sub>2</sub>N), 3.86 (2H, dm, <sup>3</sup>J<sub>HH</sub>=62Hz, <sup>3</sup>J<sub>HH</sub>=8Hz, CH<sub>2</sub>N), 2.58 (2H, dm, <sup>3</sup>J<sub>HH</sub>=40Hz, <sup>3</sup>J<sub>HH</sub>=8Hz, CH<sub>2</sub>CO), 1.27 (3H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, CH<sub>3</sub>CH<sub>2</sub>O). <sup>13</sup>C NMR (CDCl<sub>3</sub>) 171.2, 168.0, 165.6, 139.9, 139.3, 138.8, 138.1, 137.5, 132.0, 129.1, 128.5, 128.4, 128.0, 128.0, 128.0, 127.9, 127.6, 125.4, 90.6, 67.2, 60.8, 53.1, 45.0, 32.6, 14.2.

- Using the above procedure, but substituting the appropriate N-(2-amino-5-iodobenzoyl)-b-alanine ethyl ester for N-(2-diphenylmethylamino-5-iodobenzoyl)-b-alanine ethyl ester there may be prepared, for example, the following compounds:
- 1-(1-naphthyl)methyl-4-(2-carboxyethyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-(2-naphthyl)methyl-4-(2-carboxyethyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-(*m*-phenoxybenzyl)-4-(2-carboxyethyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-(*p*-phenoxybenzyl)-4-(2-carboxyethyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-(*p*-trifluoromethylbenzyl)-4-(2-carboxyethyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,

- 1-(ethyl-5-valeroyl)-4-(2-carboxyethyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,  
 1-(*p*-methoxybenzyl)-4-(2-carboxyethyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 5 1-(*p*-methylmethylbenzoyl)-4-(2-carboxyethyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,  
 1-(2-naphthyl)methyl-4-(3(*S*)-butanoate)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester,  
 1-(2-naphthyl)methyl-4-(3(*R*)-butanoate)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester,
- 10 1-(2-naphthyl)methyl-4-((+/-)-3-butanoate)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester,  
 1-(2-naphthyl)methyl-4-(3(*S*)-phenyl-3-propionate)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester,
- 15 1-(2-naphthyl)methyl-4-(3(*R*)-phenyl-3-propionate)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester,  
 1-(2-naphthyl)methyl-4-((+/-)-3-phenyl-3-propionate)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester.
- 20 c) To a magnetically stirred solution of 208 mgs of 1-(diphenylmethyl)-4-(2-carboxyethyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester (0.37mmol) in 6mL ethyl acetate, degassed of oxygen, under an atmosphere of nitrogen, was added 134 mgs of N-boc-5-amino-1-pentyne (0.74mmol), prepared as described in part (e) of Example 1, 10 mgs of
- 25 bistrisphenylphosphine palladium dichloride (0.014mmol), 5 mgs of cuprous iodide (0.026mmol) and 250uL of triethylamine (1.75mmol). After 1.5 hours, the reaction mixture was diluted with 35mL of ethyl acetate and washed with 1x15mL 5% EDTA·Na<sub>2</sub>, dried over sodium sulfate, decanted, concentrated *in vacuo*, and further purified by column
- 30 chromatography, using silica gel, eluting with a 1:1 mixture of ethyl acetate and hexane (TLC, SiO<sub>2</sub>, 1:1 EtOAc/hexane, R<sub>f</sub> = 0.46, uv positive) to yield 131 mgs (60%) of 1-(diphenylmethyl)-4-(2-carboxyethyl)-8-(N-boc-5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester. <sup>1</sup>H NMR (CDCl<sub>3</sub>, TMS) 7.71 (1H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, ArH o-CON), 7.10-7.36 (11H, m, <sup>3</sup>J<sub>HH</sub>=2Hz, ArH Ph, ArH p-CON), 6.94 (1H, d, <sup>4</sup>J<sub>HH</sub>=9Hz, ArH p-CON), 6.67 (1H, s, CHAr<sub>2</sub>), 4.78 (1H, s, NHBoc), 4.10 (2H, q, <sup>3</sup>J<sub>HH</sub>=7Hz, CH<sub>2</sub>O) 3.95 (2H, dd, <sup>2</sup>J<sub>HH</sub>=106Hz, <sup>2</sup>J<sub>HH</sub>=15Hz, COCH<sub>2</sub>CN), 3.82 (2H, dm, <sup>2,3</sup>J<sub>HH</sub>=7Hz, <sup>2,3</sup>J<sub>HH</sub>=59Hz,
- 35 CH<sub>2</sub>N), 3.19 (2H, q, <sup>3</sup>J<sub>HH</sub>=3Hz, CH<sub>2</sub>NHBoc), 2.54 (2H, dm, <sup>3</sup>J<sub>HH</sub>=7Hz, <sup>3</sup>J<sub>HH</sub>=21Hz, CH<sub>2</sub>CO), 2.37 (2H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, CH<sub>2</sub>CC), 1.71 (2H, m, <sup>3</sup>J<sub>HH</sub>=7Hz, CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>), 1.38 (9H, s, CH<sub>3</sub> Boc), 1.22 (3H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, CH<sub>3</sub>CH<sub>2</sub>O).

Using the above procedure, but substituting the appropriate 7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione for 1-(diphenylmethyl)-4-(2-carboxyethyl)-7-iodo-3,4-

dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester and alkyne for N-boc-5-amino-1-pentyne there may be prepared, for example, the following compounds:

- 1-(diphenylmethyl)-4-(2-carboxyethyl)-7-(N-boc-6-amino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 5 1-(diphenylmethyl)-4-(2-carboxyethyl)-7-(N-boc-4-amino-1-butylnl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-(1-naphthyl)methyl-4-(2-carboxyethyl)-7-(N-boc-5-amino-1-pentylnl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-(1-naphthyl)methyl-4-(2-carboxyethyl)-7-(N-boc-6-amino-1-hexynl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 10 1-(2-naphthyl)methyl-4-(2-carboxyethyl)-7-(N-boc-5-amino-1-pentylnl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-(2-naphthyl)methyl-4-(2-carboxyethyl)-7-(N-boc-6-amino-1-hexynl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 15 1-(*p*-phenoxybenzyl)-4-(2-carboxyethyl)-7-(N-boc-5-amino-1-pentylnl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-(*p*-phenoxybenzyl)-4-(2-carboxyethyl)-7-(N-boc-6-amino-1-hexynl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-(*m*-phenoxybenzyl)-4-(2-carboxyethyl)-7-(N-boc-5-amino-1-pentylnl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 20 1-(*m*-phenoxybenzyl)-4-(2-carboxyethyl)-7-(N-boc-6-amino-1-hexynl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-(*p*-trifluoromethylbenzyl)-4-(2-carboxyethyl)-7-(N-boc-5-amino-1-pentylnl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 25 1-(*p*-trifluoromethylbenzyl)-4-(2-carboxyethyl)-7-(N-boc-6-amino-1-hexynl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-(ethyl-5-valeroyl)-4-(2-carboxyethyl)-7-(N-boc-5-amino-1-pentylnl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-(ethyl-5-valeroyl)-4-(2-carboxyethyl)-7-(N-boc-6-amino-1-hexynl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 30 1-(*p*-methoxybenzyl)-4-(2-carboxyethyl)-7-(N-boc-5-amino-1-pentylnl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-(*p*-methoxybenzyl)-4-(2-carboxyethyl)-7-(N-boc-6-amino-1-hexynl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 35 1-(*p*-methylmethylbenzoyl)-4-(2-carboxyethyl)-7-(N-boc-5-amino-1-pentylnl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-(*p*-methylmethylbenzoyl)-4-(2-carboxyethyl)-7-(N-boc-6-amino-1-hexynl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,

- 1-(2-naphthyl)methyl-4-(3(S)-butanoate)-7-(N-boc-5-amino-1-pentylnl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester,  
 1-(2-naphthyl)methyl-4-(3(S)-butanoate)-7-(N-boc-6-amino-1-hexynl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester,  
 5 1-(2-naphthyl)methyl-4-(3(R)-butanoate)-7-(N-boc-5-amino-1-pentylnl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester,  
 1-(2-naphthyl)methyl-4-(3(R)-butanoate)-7-(N-boc-6-amino-1-hexynl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester,  
 1-(2-naphthyl)methyl-4-((+/-)-3-butanoate)-7-(N-boc-5-amino-1-pentylnl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester,  
 10 1-(2-naphthyl)methyl-4-((+/-)-3-butanoate)-7-(N-boc-6-amino-1-hexynl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester,  
 1-(2-naphthyl)methyl-4-(3(S)-phenyl-3-propionate)-7-(N-boc-5-amino-1-pentylnl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester,  
 15 1-(2-naphthyl)methyl-4-(3(S)-phenyl-3-propionate)-7-(N-boc-6-amino-1-hexynl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester,  
 1-(2-naphthyl)methyl-4-(3(R)-phenyl-3-propionate)-7-(N-boc-5-amino-1-pentylnl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester,  
 1-(2-naphthyl)methyl-4-(3(R)-phenyl-3-propionate)-7-(N-boc-6-amino-1-hexynl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester,  
 20 1-(2-naphthyl)methyl-4-((+/-)-3-phenyl-3-propionate)-7-(N-boc-5-amino-1-pentylnl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester,  
 1-(2-naphthyl)methyl-4-((+/-)-3-phenyl-3-propionate)-7-(N-boc-6-amino-1-hexynl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester.

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- d) To a magnetically stirred solution of 66 mgs of 1-(diphenylmethyl)-4-(2-carboxyethyl)-8-(N-boc-5-amino-1-pentylnl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester (0.11mmol) in 2mL ethyl acetate, covered with an atmosphere of hydrogen, was added 23 mgs of palladium/carbon catalyst. The mixture was stirred for 1 hr. before filtering the mixture through Celite® and concentrating *in vacuo* to yield quantitative yield of 1-(diphenylmethyl)-4-(2-carboxyethyl)-8-(N-boc-5-aminopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester. <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 7.54 (1H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, Ar-H o-CON), 6.90-7.38 (12H, m, <sup>3</sup>J<sub>HH</sub>=2Hz, Ar-H Ph, ArH p-CON, ArH m-CON), 6.54 (1H, s, NCHAr<sub>2</sub>), 4.64 (1H, s, NHBoc), 4.12 (2H, q, <sup>3</sup>J<sub>HH</sub>=7Hz, CH<sub>2</sub>O) 3.97 (2H, dd, <sup>2</sup>J<sub>HH</sub>=129Hz, <sup>2</sup>J<sub>HH</sub>=15Hz, COCH<sub>2</sub>CN), 3.82 (2H, dm, <sup>2,3</sup>J<sub>HH</sub>=6Hz, <sup>2,3</sup>J<sub>HH</sub>=75Hz, CH<sub>2</sub>N), 3.07 (2H, q, <sup>3</sup>J<sub>HH</sub>=3Hz, CH<sub>2</sub>NHBoc), 2.55 (4H, dm, t, <sup>3</sup>J<sub>HH</sub>=7Hz, <sup>3</sup>J<sub>HH</sub>=30Hz, <sup>3</sup>J<sub>HH</sub>=7Hz, CH<sub>2</sub>CO, CH<sub>2</sub>Ar), 1.3-1.6 (4H, m, m, <sup>3</sup>J<sub>HH</sub>=7Hz, <sup>3</sup>J<sub>HH</sub>=7Hz, CH<sub>2</sub>CH<sub>2</sub>Ar, CH<sub>2</sub>CH<sub>2</sub>NHBoc), 1.1-1.3 (14H, s, m, t, <sup>3</sup>J<sub>HH</sub>=7Hz, <sup>3</sup>J<sub>HH</sub>=7Hz, CH<sub>3</sub> Boc, CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>NHBoc, CH<sub>3</sub>CH<sub>2</sub>).

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Using the above procedure, but substituting the appropriate 4-(2-carboxyethyl)-7-(N-boc-amino-1-alkynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester for 1-(diphenylmethyl)-4-(2-carboxyethyl)-7-(N-boc-5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester there may be prepared, for example, the following

- 5 compounds:
- 1-(diphenylmethyl)-4-(2-carboxyethyl)-7-(N-boc-6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-(diphenylmethyl)-4-(2-carboxyethyl)-7-(N-boc-4-aminobutyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 10 1-(1-naphthyl)methyl-4-(2-carboxyethyl)-7-(N-boc-5-aminopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-(1-naphthyl)methyl-4-(2-carboxyethyl)-7-(N-boc-6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-(2-naphthyl)methyl-4-(2-carboxyethyl)-7-(N-boc-5-aminopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 15 1-(2-naphthyl)methyl-4-(2-carboxyethyl)-7-(N-boc-6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-(*p*-phenoxybenzyl)-4-(2-carboxyethyl)-7-(N-boc-5-aminopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 20 1-(*p*-phenoxybenzyl)-4-(2-carboxyethyl)-7-(N-boc-6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-(*m*-phenoxybenzyl)-4-(2-carboxyethyl)-7-(N-boc-5-aminopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-(*m*-phenoxybenzyl)-4-(2-carboxyethyl)-7-(N-boc-6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 25 1-(*p*-trifluoromethylbenzyl)-4-(2-carboxyethyl)-7-(N-boc-5-aminopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-(*p*-trifluoromethylbenzyl)-4-(2-carboxyethyl)-7-(N-boc-6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 30 1-(ethyl-5-valeroyl)-4-(2-carboxyethyl)-7-(N-boc-5-aminopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-(ethyl-5-valeroyl)-4-(2-carboxyethyl)-7-(N-boc-6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-(*p*-methoxybenzyl)-4-(2-carboxyethyl)-7-(N-boc-5-aminopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 35 1-(*p*-methoxybenzyl)-4-(2-carboxyethyl)-7-(N-boc-6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-(*p*-methylmethylbenzoyl)-4-(2-carboxyethyl)-7-(N-boc-5-aminopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,

- 1-(*p*-methylmethylbenzoyl)-4-(2-carboxyethyl)-7-(N-boc-6-amino hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,  
 1-(2-naphthyl)methyl-4-(3(S)-butanoate)-7-(N-boc-5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester,  
 5 1-(2-naphthyl)methyl-4-(3(S)-butanoate)-7-(N-boc-6-amino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester,  
 1-(2-naphthyl)methyl-4-(3(R)-butanoate)-7-(N-boc-5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester,  
 1-(2-naphthyl)methyl-4-(3(R)-butanoate)-7-(N-boc-6-amino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester,  
 10 1-(2-naphthyl)methyl-4-((+/-)-3-butanoate)-7-(N-boc-5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester,  
 1-(2-naphthyl)methyl-4-((+/-)-3-butanoate)-7-(N-boc-6-amino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester,  
 15 1-(2-naphthyl)methyl-4-(3(S)-phenyl-3-propionate)-7-(N-boc-5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester,  
 1-(2-naphthyl)methyl-4-(3(S)-phenyl-3-propionate)-7-(N-boc-6-amino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester,  
 1-(2-naphthyl)methyl-4-(3(R)-phenyl-3-propionate)-7-(N-boc-5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester,  
 20 1-(2-naphthyl)methyl-4-(3(R)-phenyl-3-propionate)-7-(N-boc-6-amino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester,  
 1-(2-naphthyl)methyl-4-((+/-)-3-phenyl-3-propionate)-7-(N-boc-5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester,  
 25 1-(2-naphthyl)methyl-4-((+/-)-3-phenyl-3-propionate)-7-(N-boc-6-amino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester.

- e) To a solution of 33 mgs of 1-(diphenylmethyl)-4-(2-carboxyethyl)-7-(N-boc-5-aminopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester (0.05mmol) in 2mL  
 30 ethyl acetate was added 0.4mL triethylsilane (2.5mmol) and 6mL of saturated HCl/ethyl acetate at 5°C. After 1 hour at 5°C the mixture was concentrated *in vacuo* and diluted with 2mL of methanol. To the methanolic solution was added 2mL of 2N sodium hydroxide. After 1 hour, the reaction was quenched with 1/2mL of acetic acid, concentrated *in vacuo*, diluted with 5mL of methanol and water and purified by high pressure liquid chromatography, using  
 35 a 1/2" C-18 reverse phase column, eluting with a solvent gradient of 30:70 methanol (0.1% trifluoroacetic acid)/water (0.1% trifluoroacetic acid), time 0 to 10 minutes, to 70:30 methanol (0.1% trifluoroacetic acid)/water (0.1% trifluoroacetic acid), time 10 minutes to 40 minutes, flow = 10mL/min. ( $R_t$  = 37.5 min., uv detection 254nm) to yield 14 mgs (46%) of 1-(diphenylmethyl)-4-(2-carboxyethyl)-7-(5-aminopentyl)-3,4-dihydro-1H-1,4-



benzodiazepine-2,5-dione trifluoroacetate. HRMS (FAB) molecular ion  $m/z=500.2529$  (calc.  $C_{30}N_3O_4H_{33}$ , 500.2549).  $^1H$  NMR ( $CDCl_3$ , TMS) 6.60-7.20 (13H, m,  $^3J_{HH}=2Hz$ , ArH Ph, ArH *p*-CON, ArH *o*-CON, ArH *m*-CON), 6.33 (1H, s, NCH-Ar<sub>2</sub>), 3.65 (2H, dd,  $^2J_{HH}=113Hz$ ,  $^2J_{HH}=15Hz$ , COCH<sub>2</sub>CN), 3.50 (2H, dm,  $^2,^3J_{HH}=7Hz$ ,  $^2,^3J_{HH}=59Hz$ , CH<sub>2</sub>N), 2.66 (2H, t,  $^3J_{HH}=8Hz$ , CH<sub>2</sub>NH<sub>2</sub>) 2.22-2.39 (4H, m, t,  $^3J_{HH}=6Hz$ ,  $^3J_{HH}=8Hz$ , CH<sub>2</sub>CO, CH<sub>2</sub>Ar), 1.20-1.33 (4H, m, m,  $^3J_{HH}=7Hz$ , CH<sub>2</sub>CH<sub>2</sub>Ar, CH<sub>2</sub>CH<sub>2</sub>NH<sub>2</sub>), 1.00 (2H, m,  $^3J_{HH}=7Hz$ , CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>NH<sub>2</sub>).

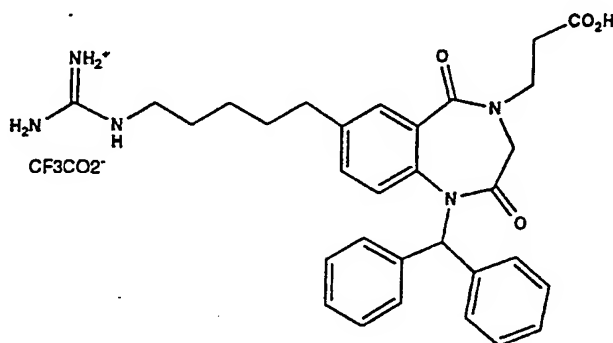
- Using the above procedure, but substituting the appropriate 3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione for 1-(diphenylmethyl)-4-(2-carboxyethyl)-7-(N-boc-5-aminopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester there may be prepared, for example, the following compounds:
- 1-(diphenylmethyl)-4-(2-carboxyethyl)-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(diphenylmethyl)-4-(2-carboxyethyl)-7-(4-aminobutyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(1-naphthyl)methyl-4-(2-carboxyethyl)-7-(5-aminopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(1-naphthyl)methyl-4-(2-carboxyethyl)-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(2-naphthyl)methyl-4-(2-carboxyethyl)-7-(5-aminopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(2-naphthyl)methyl-4-(2-carboxyethyl)-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(*p*-phenoxybenzyl)-4-(2-carboxyethyl)-7-(5-aminopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(*p*-phenoxybenzyl)-4-(2-carboxyethyl)-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(*m*-phenoxybenzyl)-4-(2-carboxyethyl)-7-(5-aminopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(*m*-phenoxybenzyl)-4-(2-carboxyethyl)-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(*p*-trifluoromethylbenzyl)-4-(2-carboxyethyl)-7-(5-aminopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(*p*-trifluoromethylbenzyl)-4-(2-carboxyethyl)-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(ethyl-5-valeroyl)-4-(2-carboxyethyl)-7-(5-aminopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(ethyl-5-valeroyl)-4-(2-carboxyethyl)-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- 1-(*p*-methoxybenzyl)-4-(2-carboxyethyl)-7-(5-aminopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-(*p*-methoxybenzyl)-4-(2-carboxyethyl)-7-(6-aminoethyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 1-(*p*-methylmethylbenzoyl)-4-(2-carboxyethyl)-7-(5-aminopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-(*p*-methylmethylbenzoyl)-4-(2-carboxyethyl)-7-(6-aminoethyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-(diphenylmethyl)-4-(2-carboxyethyl)-7-(6-amino-1-hexynyl)-3,4-dihydro-1H-
- 10 1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-(diphenylmethyl)-4-(2-carboxyethyl)-7-(4-amino-1-butylnl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-(1-naphthyl)methyl-4-(2-carboxyethyl)-7-(5-amino-1-pentynl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 1-(1-naphthyl)methyl-4-(2-carboxyethyl)-7-(6-amino-1-hexynl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-(2-naphthyl)methyl-4-(2-carboxyethyl)-7-(5-amino-1-pentynl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-(2-naphthyl)methyl-4-(2-carboxyethyl)-7-(6-amino-1-hexynl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 1-(*p*-phenoxybenzyl)-4-(2-carboxyethyl)-7-(5-amino-1-pentynl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-(*p*-phenoxybenzyl)-4-(2-carboxyethyl)-7-(6-amino-1-hexynl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 1-(*m*-phenoxybenzyl)-4-(2-carboxyethyl)-7-(5-amino-1-pentynl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-(*m*-phenoxybenzyl)-4-(2-carboxyethyl)-7-(6-amino-1-hexynl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-(*p*-trifluoromethylbenzyl)-4-(2-carboxyethyl)-7-(5-amino-1-pentynl)-3,4-dihydro-1H-
- 30 1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-(*p*-trifluoromethylbenzyl)-4-(2-carboxyethyl)-7-(6-amino-1-hexynl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-(ethyl-5-valeroyl)-4-(2-carboxyethyl)-7-(5-amino-1-pentynl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 1-(ethyl-5-valeroyl)-4-(2-carboxyethyl)-7-(6-amino-1-hexynl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-(*p*-methoxybenzyl)-4-(2-carboxyethyl)-7-(5-amino-1-pentynl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- 1-(p-methoxybenzyl)-4-(2-carboxyethyl)-7-(6-amino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(p-methylmethylbenzoyl)-4-(2-carboxyethyl)-7-(5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 1-(p-methylmethylbenzoyl)-4-(2-carboxyethyl)-7-(6-amino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(2-naphthyl)methyl-4-(3(S)-butanoate)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(2-naphthyl)methyl-4-(3(S)-butanoate)-7-(6-amino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 1-(2-naphthyl)methyl-4-(3(R)-butanoate)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(2-naphthyl)methyl-4-(3(R)-butanoate)-7-(6-amino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 1-(2-naphthyl)methyl-4-((+/-)-3-butanoate)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(2-naphthyl)methyl-4-((+/-)-3-butanoate)-7-(6-amino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(2-naphthyl)methyl-4-(3(S)-phenyl-3-propionate)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 1-(2-naphthyl)methyl-4-(3(S)-phenyl-3-propionate)-7-(6-amino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(2-naphthyl)methyl-4-(3(R)-phenyl-3-propionate)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 1-(2-naphthyl)methyl-4-(3(R)-phenyl-3-propionate)-7-(6-amino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(2-naphthyl)methyl-4-((+/-)-3-phenyl-3-propionate)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(2-naphthyl)methyl-4-((+/-)-3-phenyl-3-propionate)-7-(6-amino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 1-(2-naphthyl)methyl-4-(3(S)-butanoate)-7-(5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(2-naphthyl)methyl-4-(3(S)-butanoate)-7-(6-amino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 1-(2-naphthyl)methyl-4-(3(R)-butanoate)-7-(5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(2-naphthyl)methyl-4-(3(R)-butanoate)-7-(6-amino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- 1-(2-naphthyl)methyl-4-((+/-)3-butanoate)-7-(5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(2-naphthyl)methyl-4-((+/-)3-butanoate)-7-(6-amino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 1-(2-naphthyl)methyl-4-(3(S)-phenyl-3-propionate)-7-(5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(2-naphthyl)methyl-4-(3(S)-phenyl-3-propionate)-7-(6-amino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(2-naphthyl)methyl-4-(3(R)-phenyl-3-propionate)-7-(5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 1-(2-naphthyl)methyl-4-(3(R)-phenyl-3-propionate)-7-(6-amino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(2-naphthyl)methyl-4-((+/-)3-phenyl-3-propionate)-7-(5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 1-(2-naphthyl)methyl-4-((+/-)3-phenyl-3-propionate)-7-(6-amino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate.

## Example 15



20

*1-(diphenylmethyl)-4-(2-carboxyethyl)-7-(5-guanidinopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate*

- 25 a) To a solution of 11 mgs of 1-(diphenylmethyl)-4-(2-carboxyethyl)-7-(5-aminopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate (0.02mmol) in 2mL of methanol was added 1mL of 5% potassium bicarbonate and 17 mgs of aminoiminomethanesulfonic acid (0.12mmol). After 30 minutes, the reaction mixture was quenched with 0.5mL of acetic acid and concentrated *in vacuo*. The resulting residue was diluted with 5mL of water and
- 30 methanol and purified by high pressure liquid chromatography, using a 1/2" C-18 reverse phase column, eluting with a solvent gradient of 30:70 methanol (0.1% trifluoroacetic acid)/water (0.1% trifluoroacetic acid), time 0 to 10 minutes, to 70:30 methanol (0.1%

trifluoroacetic acid)/water (0.1% trifluoroacetic acid), time 10 minutes to 40 minutes, flow = 10mL/min. ( $R_t$  = 40.5 min., uv detection 254nm) to yield 9.3 mg (71%) of 1-(diphenylmethyl)-

4-(2-carboxyethyl)-7-(5-guanidinopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate. HRMS (FAB) molecular ion  $m/z$ =542.2750 (calc.  $C_{31}N_5O_4H_{35}$ , 542.2767).

- 5  $^1H$  NMR ( $CDCl_3$ , dTMS) 6.78-7.22 (13H, m,  $^3J_{HH}$ =2Hz, ArH Ph, ArH p-CON, ArH o-CON, ArH m-CON), 6.45 (1H, s, NCHAR<sub>2</sub>), 3.83 (2H, dd,  $^2J_{HH}$ =113Hz,  $^2J_{HH}$ =15Hz, COCH<sub>2</sub>CN), 3.65 (2H, dm,  $^2,3J_{HH}$ =7Hz,  $^2,3J_{HH}$ =59Hz, CH<sub>2</sub>N), 2.85 (2H, t,  $^3J_{HH}$ =8Hz, CH<sub>2</sub>NH<sub>2</sub>) 2.25-2.45 (4H, t, t, CH<sub>2</sub>CO, CH<sub>2</sub>Ar), 1.31 (4H, m, m,  $^3J_{HH}$ =7Hz, CH<sub>2</sub>CH<sub>2</sub>Ar, CH<sub>2</sub>CH<sub>2</sub>NH<sub>2</sub>), 0.98 (2H, m,  $^3J_{HH}$ =7Hz, CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>NH<sub>2</sub>).

10

Using the above procedure, but substituting the appropriate amino acid for 1-(diphenylmethyl)-4-(2-carboxyethyl)-7-(5-aminopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate there may be prepared, for example, the following compounds:

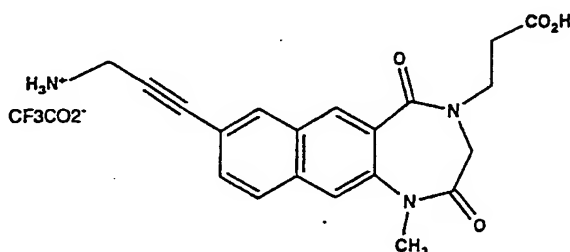
- 15 1-(diphenylmethyl)-4-(2-carboxyethyl)-7-(6-guanidinoethyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-(diphenylmethyl)-4-(2-carboxyethyl)-7-(4-guanidinobutyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-(1-naphthyl)methyl-4-(2-carboxyethyl)-7-(5-guanidinopentyl)-3,4-dihydro-1H-1,4-  
20 benzodiazepine-2,5-dione trifluoroacetate,  
1-(1-naphthyl)methyl-4-(2-carboxyethyl)-7-(6-guanidinoethyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-(2-naphthyl)methyl-4-(2-carboxyethyl)-7-(5-guanidinopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
25 1-(2-naphthyl)methyl-4-(2-carboxyethyl)-7-(6-guanidinoethyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-(p-phenoxybenzyl)-4-(2-carboxyethyl)-7-(5-guanidinopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-(p-phenoxybenzyl)-4-(2-carboxyethyl)-7-(6-guanidinoethyl)-3,4-dihydro-1H-1,4-  
30 benzodiazepine-2,5-dione trifluoroacetate,  
1-(m-phenoxybenzyl)-4-(2-carboxyethyl)-7-(5-guanidinopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-(m-phenoxybenzyl)-4-(2-carboxyethyl)-7-(6-guanidinoethyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
35 1-(p-trifluoromethylbenzyl)-4-(2-carboxyethyl)-7-(5-guanidinopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
1-(p-trifluoromethylbenzyl)-4-(2-carboxyethyl)-7-(6-guanidinoethyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- 1-(ethyl-5-valeroyl)-4-(2-carboxyethyl)-7-(5-guanidinopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(ethyl-5-valeroyl)-4-(2-carboxyethyl)-7-(6-guanidinohexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 1-(*p*-methoxybenzyl)-4-(2-carboxyethyl)-7-(5-guanidinopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(*p*-methoxybenzyl)-4-(2-carboxyethyl)-7-(6-guanidinohexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(*p*-methylmethylbenzoyl)-4-(2-carboxyethyl)-7-(5-guanidinopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 1-(*p*-methylmethylbenzoyl)-4-(2-carboxyethyl)-7-(6-guanidinohexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(diphenylmethyl)-4-(2-carboxyethyl)-7-(6-guanidino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 1-(diphenylmethyl)-4-(2-carboxyethyl)-7-(4-guanidino-1-butylnl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(1-naphthyl)methyl-4-(2-carboxyethyl)-7-(5-guanidino-1-pentylnl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(1-naphthyl)methyl-4-(2-carboxyethyl)-7-(6-guanidino-1-hexynl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 1-(2-naphthyl)methyl-4-(2-carboxyethyl)-7-(5-guanidino-1-pentylnl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(2-naphthyl)methyl-4-(2-carboxyethyl)-7-(6-guanidino-1-hexynl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 1-(*p*-phenoxybenzyl)-4-(2-carboxyethyl)-7-(5-guanidino-1-pentylnl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(*p*-phenoxybenzyl)-4-(2-carboxyethyl)-7-(6-guanidino-1-hexynl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(*m*-phenoxybenzyl)-4-(2-carboxyethyl)-7-(5-guanidino-1-pentylnl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 1-(*m*-phenoxybenzyl)-4-(2-carboxyethyl)-7-(6-guanidino-1-hexynl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(*p*-trifluoromethylbenzyl)-4-(2-carboxyethyl)-7-(5-guanidino-1-pentylnl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 1-(*p*-trifluoromethylbenzyl)-4-(2-carboxyethyl)-7-(6-guanidino-1-hexynl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(ethyl-5-valeroyl)-4-(2-carboxyethyl)-7-(5-guanidino-1-pentylnl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- 1-(ethyl-5-valeroyl)-4-(2-carboxyethyl)-7-(6-guanidino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(p-methoxybenzyl)-4-(2-carboxyethyl)-7-(5-guanidino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 5 1-(p-methoxybenzyl)-4-(2-carboxyethyl)-7-(6-guanidino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(p-methylmethylbenzoyl)-4-(2-carboxyethyl)-7-(5-guanidino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(p-methylmethylbenzoyl)-4-(2-carboxyethyl)-7-(6-guanidino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 1-(2-naphthyl)methyl-4-(3(S)-butanoate)-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(2-naphthyl)methyl-4-(3(S)-butanoate)-7-(6-guanidino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 1-(2-naphthyl)methyl-4-(3(R)-butanoate)-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(2-naphthyl)methyl-4-(3(R)-butanoate)-7-(6-guanidino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(2-naphthyl)methyl-4-((+/-)-3-butanoate)-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 20 1-(2-naphthyl)methyl-4-((+/-)-3-butanoate)-7-(6-guanidino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(2-naphthyl)methyl-4-(3(S)-phenyl-3-propionate)-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 25 1-(2-naphthyl)methyl-4-(3(S)-phenyl-3-propionate)-7-(6-guanidino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(2-naphthyl)methyl-4-(3(R)-phenyl-3-propionate)-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(2-naphthyl)methyl-4-(3(R)-phenyl-3-propionate)-7-(6-guanidino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 30 1-(2-naphthyl)methyl-4-((+/-)-3-phenyl-3-propionate)-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(2-naphthyl)methyl-4-((+/-)-3-phenyl-3-propionate)-7-(6-guanidino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 35 1-(2-naphthyl)methyl-4-(3(S)-butanoate)-7-(5-guanidino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-(2-naphthyl)methyl-4-(3(S)-butanoate)-7-(6-guanidino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

- 1-(2-naphthyl)methyl-4-(3(R)-butanoate)-7-(5-guanidino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 1-(2-naphthyl)methyl-4-(3(R)-butanoate)-7-(6-guanidino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 5 1-(2-naphthyl)methyl-4-((+/-)3-butanoate)-7-(5-guanidino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 1-(2-naphthyl)methyl-4-((+/-)3-butanoate)-7-(6-guanidino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 1-(2-naphthyl)methyl-4-(3(S)-phenyl-3-propionate)-7-(5-guanidino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 10 1-(2-naphthyl)methyl-4-(3(S)-phenyl-3-propionate)-7-(6-guanidino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 1-(2-naphthyl)methyl-4-(3(R)-phenyl-3-propionate)-7-(5-guanidino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 15 1-(2-naphthyl)methyl-4-(3(R)-phenyl-3-propionate)-7-(6-guanidino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 1-(2-naphthyl)methyl-4-((+/-)3-phenyl-3-propionate)-7-(5-guanidino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,  
 1-(2-naphthyl)methyl-4-((+/-)3-phenyl-3-propionate)-7-(6-guanidino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate.  
 20

## Example 16



*1-methyl-4-(2-carboxyethyl)-8-(3-amino-1-propynyl)-3,4-dihydro  
 1H-1,4-naphthodiazepine-2,4-dione trifluoroacetate*

- a) [4,5]-Benzoisatoic anhydride was prepared according to the method described in part (a)  
 30 of Example 1. Thus, 35g of 3-amino-2-naphthoic acid (187mMol) was dissolved in 500mL of  
 water containing 20g of potassium carbonate. A solution of phosgene in toluene (1.93M, 290mL)  
 was added slowly with vigorous stirring. A precipitate appeared along with evolution of  
 gas. Stirring was continued for 30 min after complete addition of the phosgene. The  
 precipitate was collected by suction filtration. The solid material was washed thoroughly



with water and hexane and air dried yielding 37.7g (95%) of [4,5]-benzoisatoic anhydride. M.p. >400°C. <sup>1</sup>H NMR (CDCl<sub>3</sub>): 12.0, s, 1H; 8.95, s, 1H; 8.3, d, J=7Hz, 1H; 8.15, d, J=7Hz, 1H; 7.85, t, J=7Hz, 1H; 7.7, t, J=7Hz, 1H; 7.7, s, 1H.

- 5 b) 10.3g (48.3mMol) of [4,5]-benzoisatoic anhydride was dissolved in 50mL of anhydrous N,N-dimethylformamide (DMF) and added dropwise to an ice-cooled suspension of sodium hydride (97%, 1.25g) in 50mL of DMF with gas evolution. After stirring an additional 30min, 3.16mL (105mol%) of methyl iodide was added to the solution. After one hour, the reaction mixture was poured over ice and the precipitated product was collected by suction filtration and air dried, yielding 9.3g (85%) of N-methyl-[4,5]-benzoisatoic anhydride, a tan solid. 10 M.p. 219-222°C. <sup>1</sup>H NMR (d<sup>6</sup>-DMSO): 8.8, s, 1H; 8.15, d, J=7Hz, 1H; 8.0, d, J=7Hz, 1H; 7.85, s, 1H; 7.7, t, J=7Hz, 1H; 7.55, t, J=7Hz, 1H; 3.35, s, 3H.

Using the above procedure, but substituting the appropriate alkyl halide for methyl iodide there may be prepared, for example, the following compounds:

- 15 N-benzyl-[4,5]-benzoisatoic anhydride, N-(2,4-difluorobenzyl)-[4,5]-benzoisatoic anhydride, N-(b-naphthyl)methyl-[4,5]-benzoisatoic anhydride.

- c) A solution of 20.0g (89mMol) of N-methyl-[4,5]-benzoisatoic anhydride, 20.5g (133.5mMol) of the hydrochloride salt of ethyl 3-aminopropionate, and 18.6mL (133.5mMol) of triethylamine in 100mL of DMF was stirred at 80°C for four hr. The reaction mixture was then poured over ice containing citric acid and the precipitated product was collected by suction filtration. The crude product was recrystallized from benzene/hexane, yielding 24g (60%) of 3-(N-methylamino)-2-[N-(2-carboxyethyl)]-naphthocarboxamide ethyl ester, a yellow crystalline solid. MS (FAB) 300.1. <sup>1</sup>H NMR (CDCl<sub>3</sub>): 7.8, s, 1H; 7.65, d, J=7Hz, 1H; 20 7.6, d, J=7Hz, 1H; 7.4, t, J=7Hz, 1H; 7.2, t, J=7Hz, 1H; 6.9, br s, 1H; 6.8, s, 1H; 4.2, q, J=7Hz, 2H; 3.75, q, J=5Hz, 2H; 2.95, s, 3H; 2.65, t, J=5Hz, 2H; 1.3, t, J=7Hz, 3H. 25

Using the above procedure, but substituting the appropriate 3-aminopropionate for ethyl 3-aminopropionate there may be prepared, for example, the following compounds:

- 30 (±)-3-(N-methylamino)-2-[N-(3-butanoate)]-naphthocarboxamide ethyl ester,  
(±)-3-(N-methylamino)-2-[N-(3-phenyl-3-propionate)]-naphthocarboxamide ethyl ester,  
(±)-3-(N-benzylamino)-2-[N-(3-butanoate)]-naphthocarboxamide ethyl ester,  
(±)-3-(N-benzylamino)-2-[N-(3-phenyl-3-propionate)]-naphthocarboxamide ethyl ester,  
(±)-3-(N-2,4-difluorobenzylamino)-2-[N-(3-butanoate)]-naphthocarboxamide ethyl ester,  
35 (±)-3-(N-2,4-difluorobenzylamino)-2-[N-(3-phenyl-3-propionate)]-naphthocarboxamide ethyl ester,  
(±)-3-(N-(b-naphthyl)methylamino)-2-[N-(3-butanoate)]-naphthocarboxamide ethyl ester,  
(±)-3-(N-(b-naphthyl)methylamino)-2-[N-(3-phenyl-3-propionate)]-naphthocarboxamide ethyl ester.

d) To an ice-cooled solution of 5.0g (17.5mMol) of 3-(N-methyl)amino-2-[N-(2-carboxyethyl)]-naphthocarboxamide ethyl ester and 2.68mL (19mMol) of triethylamine in 50mL of methylene chloride was added dropwise a solution of 1.59mL 19(mMol) of bromoacetyl bromide. After warming to room temperature, the solution was washed with aqueous citric acid and aqueous sodium bicarbonate solution, dried over sodium sulfate and evaporated. This material was dissolved in DMF (20mL) and added dropwise to a suspension of 97% sodium hydride (462mg, 19mMol) at -15°C. After warming to room temperature the solution was poured over ice/citric acid and extracted into ether. 3.1g (52%) of 1-methyl-4-(carboxyethyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione ethyl ester, a yellow oil, was obtained after column chromatography on silica gel. MS (FAB) 341. <sup>1</sup>H NMR (CDCl<sub>3</sub>): 8.4, s, 1H; 7.9, d, J=8.5, 1H; 7.8, d, J=8.5, 1H; 7.6, s, 1H; 7.45-7.6, m, 2H; 4.15, q, J=7Hz, 2H; 4.1, d, J=14Hz, 1H; 3.95, t, J=7Hz, 2H; 3.85, d, J=14Hz, 1H; 3.5, s, 3H; 2.75, m, 2H; 1.25, t, J=7Hz, 3H.

Using the above procedure, but substituting the appropriate naphthocarboxamide for 3-(N-methyl)amino-2-[N-(2-carboxyethyl)]-naphthocarboxamide ethyl ester there may be prepared, for example, the following compounds:

- 1-benzyl-4-(carboxyethyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione ethyl ester
- 1-(2,4-difluorobenzyl)-4-(carboxyethyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione ethyl ester
- 1-(b-naphthyl)methyl-4-(carboxyethyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione ethyl ester
- (±)-1-methyl-4-(3-butanoate)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione ethyl ester,
- (±)-1-methyl-4-(3-phenyl-3-propionate)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione ethyl ester,
- (±)-1-benzyl-2-(3-butanoate)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione ethyl ester,
- (±)-1-benzyl-2-(3-phenyl-3-propionate)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione ethyl ester,
- (±)-1-(2,4-difluorobenzyl)-2-(3-butanoate)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione ethyl ester,
- (±)-1-(2,4-difluorobenzyl)-2-(3-phenyl-3-propionate)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione ethyl ester,
- (±)-1-(b-naphthyl)methyl-2-(3-butanoate)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione ethyl ester,
- (±)-1-(b-naphthyl)methyl-4-(3-phenyl-3-propionate)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione ethyl ester.

e) 1-Methyl-4-(ethoxycarbonyl)ethyl-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione (3.1g, 9.1mMol) was placed in an ice bath and 4mL of fuming nitric acid was added. After

stirring for 1hr at 0°C the reaction was allowed to warm to room temperature and stirred for an additional 2hr. The reaction mixture was poured into ice water and extracted with ethyl acetate, dried (sodium sulfate) and evaporated to yield a yellow solid. Thin layer chromatography (TLC) and <sup>1</sup>H NMR analysis indicated a mixture of nitrated products. The crude mixture was purified by column chromatography (SiO<sub>2</sub>) to yield 600mgs (17%) of 8-Nitro-1-methyl-4-(ethoxycarbonylethyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione. <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 8.8 (1H, d, <sup>4</sup>J<sub>HH</sub>=2Hz), 8.6 (1H, s), 8.3 (1H, dd, <sup>2</sup>J<sub>HH</sub>=9Hz, <sup>4</sup>J<sub>HH</sub>=2Hz), 7.95 (1H, d, <sup>3</sup>J<sub>HH</sub>=9Hz), 7.7 (1H, s), 4.15 (2H, q, <sup>3</sup>J<sub>HH</sub>=8Hz), 4.1 (1H, d, <sup>2</sup>J<sub>HH</sub>=14Hz), 4.0 (2H, t, <sup>3</sup>J<sub>HH</sub>=7Hz), 3.95 (1H, <sup>2</sup>J<sub>HH</sub>=14Hz), 3.5 (3H, s), 2.75 (2H, m), 1.25 (3H, t, <sup>3</sup>J<sub>HH</sub>=8Hz).

Using the above procedure, but substituting the appropriate naphthocarboxamide for 3-(N-methyl)amino-2-[N-(2-carboxyethyl)]-naphthocarboxamide ethyl ester there may be prepared, for example, the following compounds:

1-(2,4-difluorobenzyl)-4-(carboxyethyl)-8-nitro-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione ethyl ester,  
 (±)-1-methyl-4-(3-butanoate)-8-nitro-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione ethyl ester,  
 (±)-1-(2,4-difluorobenzyl)-2-(3-butanoate)-8-nitro-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione ethyl ester.

f) 1-Methyl-4-(carboxylethyl)-8-nitro-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione ethyl ester (600mg, 1.56mMol) was hydrogenated in ethanol with 200mg of 10% palladium on carbon catalyst at 50psi hydrogen for 24hr after which time tlc indicated that reduction was complete. The catalyst was removed by filtration through Celite®, and the filtrate was evaporated yielding ca. 500mg of 1-methyl-4-(carboxylethyl)-8-amino-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione ethyl ester. <sup>1</sup>H NMR (CDCl<sub>3</sub>): 8.1, s, 1H; 7.6, d, J=9.5Hz, 1H; 7.45, s, 1H; 7.15, s, 1H; 7.1, d, J=9.5Hz, 1H; 4.15, q, J=8Hz, 2H; 4.1, d, J=14Hz, 1H; 3.95, t, J=7Hz, 2H; 3.8, d, J=14Hz, 1H; 3.4, s, 3H; 2.7, m, 2H; 1.2, t, J=8Hz. This material was dissolved in 4mL of water containing 117mL of concentrated sulfuric acid and cooled to 0°C with an ice bath. Sodium nitrite (97mg, 1.5mMol) in 2mL of water was added to the solution dropwise, and the temperature of the solution was allowed to rise to room temperature over one hour. 2.6g (14mMol) of potassium iodide was added to the solution and stirring was continued for 24hr. The reaction mixture was diluted with water and extracted with ethyl acetate, and the ethyl acetate was dried (sodium sulfate) and evaporated. Chromatography on silica gel yielded ca. 240 mg of 8-Iodo-1-methyl-4-(ethoxycarbonylethyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione as an amber oil. MS (FAB) 466.9. <sup>1</sup>H NMR (CDCl<sub>3</sub>): 8.3, s, 2H; 7.8, dd, J=8.5Hz, J'=1.5Hz, 1H; 7.55, s, 1H; 7.55, d, J=8.5Hz, 1H; 4.15, q, J=7Hz, 2H; 4.1, d, J=14Hz, 1H; 3.95, t, J=7Hz, 2H; 3.9, d, J=14Hz, 1H; 3.5, s, 3H; 2.75, m, 2H; 1.25, t, J=7Hz.

Using the above procedure, but substituting the appropriate 8-nitro-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione for 1-methyl-4-(2-carboxylethyl)-8-nitro-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione ethyl ester there may be prepared, for example, the following compounds:

- 5 1-(2,4-difluorobenzyl)-4-(carboxyethyl)-8-iodo-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione ethyl ester,  
(±)-1-methyl-4-(3-butanoate)-8-iodo-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione ethyl ester,  
(±)-1-(2,4-difluorobenzyl)-2-(3-butanoate)-8-iodo-3,4-dihydro-1H-1,4-naphthodiazepine-  
10 2,5-dione ethyl ester.

- g) 1-Methyl-4-(ethoxycarbonylethyl)-8-iodo-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione (100mg, 215mM) was combined in 1mL of ethyl acetate with 166mg (500m%) of N-t-butoxycarbonylpropargyl amine, 15mg (10mol%) of bistrisphenylphosphine palladium  
15 dichloride, and 299mL (10X excess) of triethylamine. The apparatus was degassed and an atmosphere of nitrogen was introduced. 8.2mg (20mol%) of cuprous iodide was added and the solution was again degassed and nitrogen atmosphere was replaced. After 24hr the reaction mixture was diluted with ethyl acetate and washed successively with aqueous citric acid and aqueous sodium bicarbonate, dried (sodium sulfate) and evaporated. The residue was  
20 chromatographed on silica gel, yielding 110mg of 1-methyl-4-(crboxylethyl)-8-(3-t-butyloxycarbonylamino-1-propynyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione. MS (FAB) 494.1. <sup>1</sup>H NMR (CDCl<sub>3</sub>): 8.35, s, 1H; 7.95, s, 1H; 7.75, d, J=8.5Hz, 1H; 7.55, s, 1H; 7.5, d, J=8.5Hz, 1H; 4.9, br s, 1H; 4.2, br s, 2H; 4.2, q, J=7Hz, 2H; 4.1, d, J=14Hz, 1H; 4.0, t, J=7Hz, 2H; 3.9, d, J=7Hz, 1H; 3.5, s, 3H; 2.75, m, 2H; 1.5, s, 9H; 1.25, t, J=7Hz, 3H.

- 25 Using the above procedure, but substituting the appropriate 8-iodo-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione for 1-methyl-4-(2-carboxylethyl)-8-iodo-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione ethyl ester there may be prepared, for example, the following compounds:

- 1-(2,4-difluorobenzyl)-4-(carboxyethyl)-8-(3-t-butyloxycarbonylamino-1-propynyl)-3,4-  
30 dihydro-1H-1,4-naphthodiazepine-2,5-dione ethyl ester,  
(±)-1-methyl-4-(3-butanoate)-8-(3-t-butyloxycarbonylamino-1-propynyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione ethyl ester,  
(±)-1-(2,4-difluorobenzyl)-2-(3-butanoate)-8-(3-t-butyloxycarbonylamino-1-propynyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione ethyl ester,  
35 1-(2,4-difluorobenzyl)-4-(carboxyethyl)-8-(4-t-butyloxycarbonylamino-1-butynyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione ethyl ester,  
(±)-1-methyl-4-(3-butanoate)-8-(4-t-butyloxycarbonylamino-1-butynyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione ethyl ester,

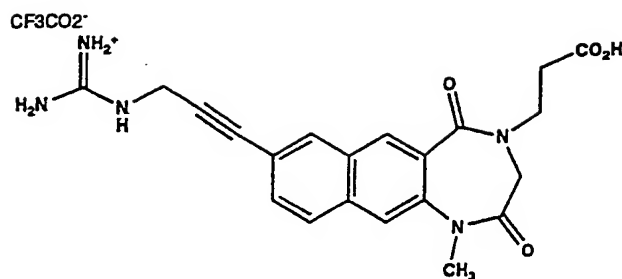
(±)-1-(2,4-difluorobenzyl)-2-(3-butanoate)-8-(4-t-butyloxycarbonylamino-1-butynyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione ethyl ester.

- h) To a stirred solution of 1-methyl-4-(carboxyethyl)-8-(3-t-butyloxycarbonylamino-1-propynyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione ethyl ester (110mg, 0.22mMol) in 8mL of methanol was added 1mL of 1N sodium hydroxide solution. After 4hr no more starting material was detected by tlc. The solvent was removed *in vacuo* and the residue was treated with 10mL of 4N HCl in dioxane. After 2hr the solvent was removed *in vacuo* and the residue was purified by reverse phase HPLC, yielding 25mg of 1-methyl-4-(carboxyethyl)-8-(3-amino-1-propynyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione. HRMS (FAB) molecular ion  $m/z=366.1454$  (calcd.  $C_{20}H_{20}N_3O_4$ , 366.4004).  $^1H$  NMR ( $CD_3OD$ ): 7.9, s, 1H; 7.7, s, 1H; 7.55, d,  $J=8Hz$ , 1H; 7.45, s, 1H; 7.2, d,  $J=8Hz$ , 1H; 3.75, d,  $J=14Hz$ , 1H; 3.65, s, 2H; 3.55, d,  $J=14Hz$ , 1H; 3.5, m, 2H; 3.05, s, 3H; 2.2, q,  $J=7Hz$ , 2H.

- 15 Using the above procedure, but substituting the appropriate 8-(t-butyloxycarbonylamino-1-alkynyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione for 1-methyl-4-(carboxyethyl)-8-(3-t-butyloxycarbonylamino-1-propynyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione ethyl ester there may be prepared, for example, the following compounds:

- 20 1-(2,4-difluorobenzyl)-4-(carboxyethyl)-8-(3-amino-1-propynyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione,  
(±)-1-methyl-4-(3-butanoate)-8-(3-amino-1-propynyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione,  
(±)-1-(2,4-difluorobenzyl)-4-(3-butanoate)-8-(3-amino-1-propynyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione,  
25 1-(2,4-difluorobenzyl)-4-(carboxyethyl)-8-(4-amino-1-butynyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione,  
(±)-1-methyl-4-(3-butanoate)-8-(4-amino-1-butynyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione,  
30 (±)-1-(2,4-difluorobenzyl)-4-(3-butanoate)-8-(4-amino-1-butynyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione.

## Example 17



5

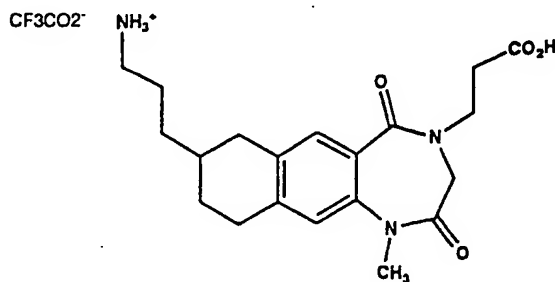
*1-Methyl-4-(carboxyethyl)-8-(3-guanidino-1-propynyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione*

The title compound was prepared from 1-methyl-4-(carboxyethyl)-8-(3-amino-1-propynyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione (50mg) according to the procedure described in Example 2. HRMS (FAB) molecular ion  $m/z=408.1672$  (calcd.  $C_{21}H_{22}N_5O_4$ , 408.1672).  $^1H$  NMR ( $CD_3OD$ ): 7.9, s, 1H; 7.7, s, 1H; 7.55, d,  $J=8Hz$ , 1H; 7.45, s, 1H; 7.2, d,  $J=8Hz$ , 1H; 3.75, d,  $J=14Hz$ , 1H; 3.65, s, 2H; 3.55, d,  $J=14Hz$ , 1H; 3.5, m, 2H; 3.05, s, 3H; 2.2, q,  $J=7Hz$ , 2H.

Using the above procedure, but substituting the appropriate 8-(amino-1-alkynyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione for 1-methyl-4-(carboxyethyl)-8-(3-amino-1-propynyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione there may be prepared, for example, the following compounds:

- 1-(2,4-difluorobenzyl)-4-(carboxyethyl)-8-(3-guanidino-1-propynyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione,
- (±)-1-methyl-4-(3-butanoate)-8-(3-guanidino-1-propynyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione,
- (±)-1-(2,4-difluorobenzyl)-2-(3-butanoate)-8-(3-guanidino-1-propynyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione,
- 1-(2,4-difluorobenzyl)-4-(carboxyethyl)-8-(4-guanidino-1-butynyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione,
- (±)-1-methyl-4-(3-butanoate)-8-(4-guanidino-1-butynyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione,
- (±)-1-(2,4-difluorobenzyl)-2-(3-butanoate)-8-(4-guanidino-1-butynyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione.

## Example 18



5     **1-Methyl-4-(carboxyethyl)-8-(3-amino-1-propyl)-3,4,7,8,9,10-hexahydro-1H-1,4-naphthodiazepine-2,5-dione**

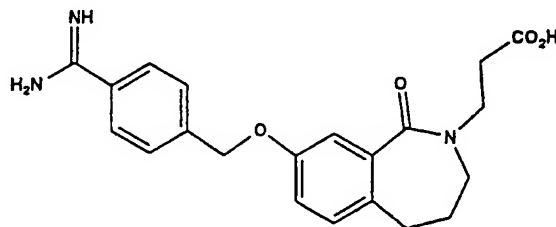
1-Methyl-4-(carboxyethyl)-8-(3-amino-1-propynyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione (50mg) was reduced with 10mg of 10% Pd/C in 10mL of 5:1 water ethanol at 45psi of hydrogen in a Parr hydrogenation apparatus for 24hr yielding, after  
 10     filtration through Celite® and purification by reverse phase HPLC, the title compound. MS (FAB) 374.1. <sup>1</sup>H NMR (CD<sub>3</sub>OD): 7.4, s, 1H; 7.0, s, 1H; 3.95, d, J=14Hz, 1H; 3.8, d, J=14Hz, 1H; 3.6, m, 2H; 3.2, s, 3H; 2.85, t, J=7Hz, 2H; 2.4, t, J=7Hz, 2H; 1.9, m, 1H; 1.65, m, 4H; 1.4, m, 4H.

Using the above procedure, but substituting the appropriate 8-(amino-1-alkynyl)-3,4-dihydro-1H-1,4-naphthodiazepine-2,5-dione for 1-methyl-4-(carboxyethyl)-8-(3-amino-1-propynyl)-3,4,7,8,9,10-hexahydro-1H-1,4-naphthodiazepine-2,5-dione there may be prepared, for example, the following compounds:

- 1-(2,4-difluorobenzyl)-4-(carboxyethyl)-8-(3-guanidino-1-propynyl)-3,4,7,8,9,10-hexahydro-1H-1,4-naphthodiazepine-2,5-dione,  
 20     (±)-1-methyl-4-(3-butanoate)-8-(3-guanidino-1-propynyl)-3,4,7,8,9,10-hexahydro-1H-1,4-naphthodiazepine-2,5-dione,  
 (±)-1-(2,4-difluorobenzyl)-2-(3-butanoate)-8-(3-guanidino-1-propynyl)-3,4,7,8,9,10-hexahydro-1H-1,4-naphthodiazepine-2,5-dione,

## Example 19

25



**2-(2-carboxyethyl)-8-(4-amidinobenzoyloxy)-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one trifluoroacetate.**

- a) A solution of 7-methoxytetralone (Fluka Chemical, 100.0 g, 0.567 mol) in dry pyridine (600 mL) was treated with hydroxylamine hydrochloride (Fluka, 43.0 g, 0.618 mol) and the resulting dark yellow solution was stirred at room temperature for 1 hour. The solution was then concentrated *in vacuo*, and azeotroped with several volumes of toluene, then partitioned between ethyl acetate and brine. The combined organic layers were dried ( $K_2CO_3$ ), filtered, and concentrated *in vacuo*. The residue was crystallized from ethyl acetate/hexanes to yield 88 g (81%) of the oxime as a waxy solid.
- b) Following a procedure of Tomita, et. al., *J. Chem. Soc. C* 1969, 183, 7-methoxytetralone oxime (45.0 g, 0.235 mol) was mixed with solid trichloroacetic acid (150 g), and the resulting suspension heated to ca. 80 C, at which time a violent exotherm ensued, yielding a black solid mixture. The residue was cooled, partitioned between ethyl acetate and saturated aqueous sodium bicarbonate solution, and the combined organic layers were concentrated *in vacuo*. The residue was purified by chromatography over silica gel to yield 10.0 g (22%) of 8-methoxy-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one. An analytical sample was obtained by recrystallization from  $CH_2Cl_2$  /hexanes to yield colorless crystals: mp. 99-101°C (lit. (see above) 100-101°C);  $^1H$  NMR( $CDCl_3$ )  $\delta$ 7.35(brs, 1H), 7.30(d, J=3Hz, 1H), 7.10(d, J=9Hz, 1H), 6.95(dd, J=3, 9Hz, 1H), 3.82(s, 3H), 3.10(q, J=6Hz, 2H), 2.80(t, J=6Hz, 2H), 1.98(m, 2H).
- c) A solution of 8-methoxy-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one (10.0 g, 0.052 mol) in N,N-dimethylformamide (30 mL) and allylbromide (23 mL, 0.260 mL) was treated with NaH (60% by wt. dispersion in oil, 2.50 g, 0.063 mol) and the resulting grey suspension stirred at room temperature under an atmosphere of argon. Gas evolution was noted. After 15 minutes, the solution was diluted with brine and partitioned between ethyl acetate and brine. The combined organic layers were concentrated *in vacuo* and purified by chromatography over silica gel eluting with a gradient of hexane to 1:1 ethyl acetate/hexane to yield 8 g (83%) of 2-allyl-8-methoxy-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one as a colorless oil:  $^1H$  NMR( $CDCl_3$ )  $\delta$ 7.23(m, 1H), 7.03(d, J=9Hz, 1H), 6.90(dd, J=3, 9Hz, 1H), 5.91(m, 1H), 5.23(m, 2H), 4.20(d, J=6Hz, 2H), 3.82(s, 3H), 3.18(t, J=6Hz, 2H), 2.72(t, J=6Hz, 2H), 1.96(pentet, J=6Hz, 2H);  $^{13}C$  NMR( $CDCl_3$ )  $\delta$ 170.59(C=O), 158.42(NC=O), 136.88, 133.81, 129.43, 129.35, 129.33, 117.56, 117.54, 112.98, 55.26, 49.43, 45.37, 29.57, 29.22; MS(FAB)  $m/z$  = 232.1( $MH^+$ ).
- d) A solution of 2-allyl-8-methoxy-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one (8.0 g, 0.043 mol) in dry tetrahydrofuran (40 mL) was treated with 9-borabicyclononane (1.0 M in THF) and the resulting mixture was stirred for 18 hours at room temperature. The solution was then diluted with EtOH (50 mL), cooled to 0°C (ice bath), and treated with 15% aqueous NaOH (15 mL) followed by 30% aqueous  $H_2O_2$  (10 mL). The resulting mixture was stirred for 1 hour



at 0°C, then treated cautiously with saturated aqueous sodium thiosulfate (50 mL). The mixture was partitioned between ethyl acetate and brine, and the combined organic layers were concentrated *in vacuo*, and the residue was purified by chromatography over silica gel eluting with ethyl acetate to yield 6.5 g (60%) of 2-(3-hydroxypropyl)-8-methoxy-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one as a colorless oil; <sup>1</sup>H NMR(CDCl<sub>3</sub>) δ 7.19(s, 1H), 7.06(d, J=9Hz, 1H, 6.92(dd, J=3, 9Hz, 1H), 4.30(brs, 1H), 3.82(s, 3H), 3.71(t, J=6Hz, 2H), 3.61(brs, 2H), 3.18(t, J=6Hz, 2H), 2.72(t, J=6Hz, 2H), 2.02(pentet, J=6Hz, 2H), 1.79(pentet, J=6Hz, 2H); <sup>13</sup>C NMR(CDCl<sub>3</sub>) δ 172.25(CO), 158.51(NCO), 136.34, 129.53, 117.46, 112.94, 58.19, 55.42, 46.16, 43.30, 30.91, 29.45, 29.12; MS (FAB) m/z=250.1(MH<sup>+</sup>); HRMS (FAB, MH<sup>+</sup>) m/z=249.1364 (calcd C<sub>14</sub>H<sub>19</sub>NO<sub>3</sub>: 249.1364).

e) A solution of 2-(3-hydroxypropyl)-8-methoxy-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one (6.50 g, 0.026 mol) in CH<sub>2</sub>Cl<sub>2</sub> was added, at -78°C (CO<sub>2</sub>/acetone), to a solution of (COCl)<sub>2</sub> (4.50 mL, 0.052 mol) and DMSO (7.4 mL, 0.104 mol) in CH<sub>2</sub>Cl<sub>2</sub> (100 mL) at -78°C. After stirring for 15 minutes at -78°C, the mixture was treated with NEt<sub>3</sub> (30 mL, 0.208 mol), and the bath removed. The CH<sub>2</sub>Cl<sub>2</sub> solution was washed with brine, dried (MgSO<sub>4</sub>) and concentrated in vacuo. The residue was immediately dissolved in N,N-dimethylformamide (30 mL) and MeOH (30 mL), and treated, in portions, with pyridinium dichromate (30 g total). After stirring for 24 hours, the dark brown suspension was diluted with ethyl acetate and filtered, and the filtrate was partitioned between ethyl acetate and brine. The combined organic layers were washed with water, brine, dried (MgSO<sub>4</sub>) and concentrated in vacuo. The residue was purified by chromatography over silica gel eluting with a gradient of 10% ethyl acetate/hexane to 100% ethyl acetate to yield 2.5 g (35%) of 2-(2-carboxyethyl)-8-methoxy-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one methyl ester as a colorless oil; <sup>1</sup>H NMR(CDCl<sub>3</sub>) δ 7.18(d, J=3Hz, 1H), 7.02(d, J=9Hz, 1H), 6.88(dd, J=3, 9Hz, 1H), 3.83(t, J=6Hz, 2H), 3.80(s, 3H), 3.69(s, 3H), 3.23(t, J=6Hz, 2H), 2.78(t, J=6Hz, 2H), 2.67(t, J=6Hz, 2H), 1.96(m, 2H); <sup>13</sup>C NMR(CDCl<sub>3</sub>) δ 172.41(CO), 171.15(CO), 158.50, 136.80, 129.58, 129.40, 117.34, 112.86, 55.42, 51.78, 47.37, 44.13, 33.64, 29.85, 29.14; MS(FAB) m/z=278.1 (MH<sup>+</sup>); HRMS(FAB, MH<sup>+</sup>) m/z=278.1392 (calcd for C<sub>15</sub>H<sub>19</sub>NO<sub>4</sub>: 277.1313).

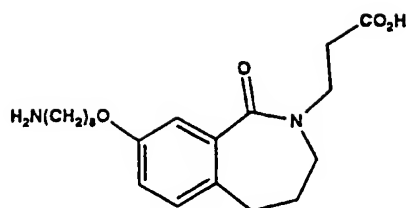
f) A solution of 2-(2-carboxyethyl)-8-methoxy-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one methyl ester (3.20 g, 0.011 mol) in CH<sub>2</sub>Cl<sub>2</sub> was added to a suspension of AlCl<sub>3</sub> (4.70 g, 0.034 mol) and *n*-propanethiol (5.3 mL, 0.057 mol) at 0°C (ice bath), and the resulting mixture stirred for 10 minutes at room temperature. The mixture was diluted with saturated aqueous sodium bicarbonate (50 mL), and the resulting white suspension stirred for 1 hour. The gelatinous mixture was filtered, and the filtrate was dried (MgSO<sub>4</sub>) and concentrated *in vacuo* to yield 2.0 g (85%) of 2-(2-carboxyethyl)-8-hydroxy-2,3,4,5-tetrahydro-1H-2-

benzazepin-1-one methyl ester as a colorless oil;  $^1\text{H}$  NMR( $\text{CD}_3\text{OD}$ )  $\delta$ 7.08(d, J=9Hz, 1H, 7.02(d, J=3Hz, 1H), 6.87(dd, J=3, 9Hz, 1H), 3.05(t, J=6Hz, 2H), 2.72(t, J=6Hz, 2H), 1.97(t, J=6Hz, 2H).

- g) A suspension of 2-(2-carboxyethyl)-8-hydroxy-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one methyl ester (250 mg, 0.950 mmol), *p*-cyanobenzylbromide (186 mg, 0.950 mmol) and  $\text{K}_2\text{CO}_3$  (500 mg) in *N,N*-dimethylformamide (20 mL) was stirred vigorously at room temperature under an atmosphere of argon. After 1 hour, the suspension was diluted with ethyl acetate (100 mL), washed with water and brine, dried ( $\text{Na}_2\text{SO}_4$ ) and concentrated *in vacuo*. The residue was purified by chromatography over silica gel to yield 150 mg (40%) of 2-(2-carboxyethyl)-8-(4-cyanobenzoyloxy)-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one methyl ester as a colorless oil;  $^1\text{H}$  NMR( $\text{CDCl}_3$ )  $\delta$ 7.66(d, J=6Hz, 1H), 7.54(d, J=6Hz, 1H), 7.00(dd, J=3, 9Hz, 1H), 5.13(s, 2H), 3.84(t, J=6Hz, 2H), 3.70(s, 3H), 3.24(t, J=6Hz, 2H), 2.72(m, 4H), 1.98(pentet, J=6Hz, 2H);  $^{13}\text{C}$  NMR( $\text{CDCl}_3$ )  $\delta$ 172.27(CO), 170.82(CO), 157.05, 142.32, 137.01, 132.39, 132.32, 130.45, 129.63, 127.52, 118.66, 117.83, 113.98, 111.53, 68.90, 51.76, 47.31, 44.14, 33.58, 29.76, 29.13; MS(FAB)  $m/z$  = 379.1( $\text{MH}^+$ ).

- h) A solution of 2-(2-carboxyethyl)-8-(4-cyanobenzoyloxy)-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one methyl ester (150 mg, 0.400 mmol) in 50 mL of 1:1 pyridine- $\text{NEt}_3$  was saturated with  $\text{H}_2\text{S}$ , and heated at  $50^\circ\text{C}$  under argon. After 1 hour, the solvents were removed *in vacuo*. The residue was dissolved in  $\text{CH}_2\text{Cl}_2$  (20 mL) and treated with iodomethane (4 mL), and the resulting mixture heated at reflux for 1 hour. The solvents were removed *in vacuo*, and the residue was dissolved in MeOH and treated with an excess of  $\text{NH}_4\text{OAc}$  (1 g). The resulting suspension was stirred for 15 minutes at  $60^\circ\text{C}$ . The MeOH was then removed, and the residue was treated with 50% NaOH (3 mL) in tetrahydrofuran (20 mL), and stirred for 15 minutes at room temperature. The mixture was then diluted with trifluoroacetic acid, and the volatiles were removed *in vacuo*. The residue was purified by HPLC (1/2" C-18 reverse-phase column, eluting with a solvent gradient of 1:9 acetonitrile(0.1%TFA)/water (0.1% TFA) time 0 to 10 minutes, to 1:1 acetonitrile(0.1%TFA)/water (0.1% TFA) time 10 to 40 minutes, flow 10mL/min,  $R_t$ =30.5 min, mu detection 254nm), to give 18.5 mg (9 %) of 2-(2-carboxyethyl)-8-(4-amidinobenzoyloxy)-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one trifluoroacetate as a yellowish-brown solid after lyophilization;  $^1\text{H}$  NMR( $\text{CD}_3\text{OD}$ )  $\delta$ 7.66(d, J=6Hz, 1H), 7.54(d, J=6Hz, 1H), 7.00(dd, J=3, 9Hz, 1H), 5.13(s, 2H), 3.84(t, J=6Hz, 2H), 3.70(s, 3H), 3.24(t, J=6Hz, 2H), 2.72(m, 4H), 1.98(m, J=6Hz, 2H); MS(FAB)  $m/z$  = 381.1688 ( $\text{MH}^+$ ).

## Example 20



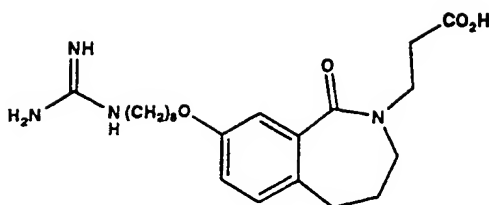
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*2-(2-carboxyethyl)-8-(8-aminoctyloxy)-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one trifluoroacetate.*

- 10 a) A solution of 2-allyl-8-methoxy-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one (1.10 g, 0.0047 mol) in  $\text{CH}_2\text{Cl}_2$  (20 mL) was treated at  $0^\circ\text{C}$  with  $\text{BBr}_3$  (1M in  $\text{CH}_2\text{Cl}_2$ , 9.50 mL), and the resulting mixture stirred at  $0^\circ\text{C}$  for 30 minutes. The solution was poured over ice,  $\text{NaHCO}_3$  was added, and the solution was extracted with  $\text{CH}_2\text{Cl}_2$ . The combined organic layers were dried ( $\text{MgSO}_4$ ), filtered, and concentrated *in vacuo* to yield ca. 1.0 g (100%) of 2-allyl-8-hydroxy-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one as a colorless oil;  $^1\text{H}$  NMR( $\text{CDCl}_3$ )
- 15 d8.2(brs, 1H), 7.47(s, 1H), 6.95(d,  $J=6\text{Hz}$ , 1H), 6.84(dd,  $J=3, 6\text{Hz}$ , 1H), 5.89(m, 1H), 5.22(m, complex, 2H), 4.20(d,  $J=6\text{Hz}$ , 2H), 3.18(t,  $J=6\text{Hz}$ , 2H), 2.66(t,  $J=6\text{Hz}$ , 2H), 2.15(s, 1H), 1.96(m, 2H);  $^{13}\text{C}$  NMR( $\text{CDCl}_3$ ) d171.65(C=O), 156.01, 135.68, 133.27, 129.55, 128.46, 118.51, 118.05, 115.91, 53.46, 49.81, 45.83, 30.87, 29.76, 29.12.
- 20 b) A suspension of 2-allyl-8-hydroxy-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one (1.0 g, 0.0049 mol), 8-azido-1-(toluenesulfonyloxy)octane (1.50 g, 0.0049 mol) and  $\text{K}_2\text{CO}_3$  in N,N-dimethylformamide (20 mL) was stirred vigorously under argon, and heated at  $65^\circ\text{C}$  for 18 hours. The solution was then partitioned between ethyl acetate and brine, and the combined organic layers were washed with water, dried ( $\text{K}_2\text{CO}_3$ ), filtered, and concentrated *in vacuo*.
- 25 The residue was purified by chromatography over silica gel to yield 920 mg (51%) of 2-allyl-8-(8-azidoctyloxy)-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one as a colorless oil;  $^1\text{H}$  NMR( $\text{CDCl}_3$ ) d7.47(s, 1H), 6.95(d,  $J=6\text{Hz}$ , 1H), 6.84(dd,  $J=3, 6\text{Hz}$ , 1H), 5.89(m, 1H), 5.22(m, complex, 2H), 4.20(d,  $J=6\text{Hz}$ , 2H), 3.95(t,  $J=6\text{Hz}$ , 2H), 3.22(t,  $J=6\text{Hz}$ , 2H), 3.18(t,  $J=6\text{Hz}$ , 2H), 2.66(t,  $J=6\text{Hz}$ , 2H), 1.90(t,  $J=6\text{Hz}$ , 2H), 1.79(t,  $J=6\text{Hz}$ , 2H), 1.60(m, 4H), 1.35(m, 6H).
- 30 c) A solution of 2-allyl-8-(8-azidoctyloxy)-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one (920 mg, 2.48 mmol) in tetrahydrofuran (20 mL) was treated with 9-borabicyclononane (0.5 M/THF, 10 mL, 5.00 mmol), and the mixture was stirred under argon at room temperature for 18 hours. The mixture was then cooled in an ice bath, and EtOH (5 mL) followed by 15%
- 35 aqueous NaOH (10 mL) and 30%  $\text{H}_2\text{O}_2$  (15 mL) were added. The mixture was stirred for 1

- hour, then saturated aqueous sodium thiosulfate was added (20 mL). The solution was partitioned between ethyl acetate and brine, and the combined ethyl acetate layers were dried (MgSO<sub>4</sub>), filtered, and concentrated *in vacuo*. The residue was purified by chromatography over silica gel to yield 400 mg (42%) of 2-(3-hydroxypropyl)-8-(8-azidooctyloxy)-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one as a colorless oil; <sup>1</sup>H NMR(CDCl<sub>3</sub>) δ 7.20(d, J=3Hz, 1H), 7.04(d, J=6Hz, 1H), 6.91(dd, J=3, 6Hz, 1H), 3.96(t, J=6Hz, 2H), 3.72(t, J=6Hz, 2H), 3.61(t, J=6Hz, 2H), 3.26(t, J=6Hz, 2H), 3.18(t, J=6Hz, 2H), 2.73(t, J=6Hz, 2H), 2.02(t, J=6Hz, 2H), 1.77(m, 4H), 1.60(m, 2H), 1.36(m, 8H).
- 10 d) 2-(3-hydroxypropyl)-8-(8-azidooctyloxy)-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one was subjected to standard Swern conditions followed by PDC oxidation as described in part (e) of Example 14 to yield 140 mg (45%) of 2-(2-carboxyethyl)-8-(8-azidooctyloxy)-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one methyl ester as a colorless oil; <sup>1</sup>H NMR(CDCl<sub>3</sub>) δ 7.18(d, J=3Hz, 1H), 7.00(d, J=9Hz, 1H), 6.89(dd, J=3, 9Hz, 1H), 3.96(t, J=6Hz, 2H), 3.84(t, J=6Hz, 2H), 3.70(s, 3H), 3.26(m, 4H), 2.74(m, 4H), 1.97(t, J=6Hz, 2H), 1.77(t, J=6Hz, 2H), 1.60(t, J=6Hz, 2H), 1.35(m, 6H); <sup>13</sup>C NMR(CDCl<sub>3</sub>) δ 172.38(C=O), 171.21(C=O), 158.04, 136.72, 129.40, 129.34, 117.78, 113.57, 68.05, 51.76, 51.41, 47.37, 44.13, 33.64, 29.86, 29.13, 19.02, 28.77, 26.59, 25.87; MS(FAB) 417.2(MH<sup>+</sup>).
- 15 20 e) A solution of 2-(2-carboxyethyl)-8-(8-azidooctyloxy)-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one methyl ester (70 mg, 0.168 mmol) in MeOH (1 mL) was treated with 10% Pd/C (20 mg) and the resulting suspension was reduced under an atmosphere of hydrogen (30psi) for 2 hours. The solution was then filtered through Celite®, and the filtrate was concentrated *in vacuo*. The residue was dissolved in tetrahydrofuran (1 mL) and treated with 50% NaOH (0.5 mL), and the resulting mixture was stirred at room temperature. After 1 hour, the solution was treated with trifluoroacetic acid (1 mL), and concentrated *in vacuo*. The white paste remaining was purified by HPLC (1/2' C-18 reverse-phase column) eluting with a gradient of 3:7 methanol (0.1% trifluoroacetate)/water (0.1% trifluoroacetate), time 0 to 10 min, to 7:3 methanol(0.1% trifluoroacetate)/water(0.1% trifluoroacetate), time 10 to 40 min, to yield 30 mg (36%) of 2-(2-carboxyethyl)-8-(8-aminooctyloxy)-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one as a colorless oil; <sup>1</sup>H NMR(CD<sub>3</sub>OD) δ 7.095(m, 2H), 6.96(dd, J=3, 9Hz, 1H), 3.98(t, J=6Hz, 2H), 3.83(t, J=6Hz, 2H), 3.27(m, 4H), 2.90(t, J=6Hz, 2H), 2.69(m, 4H), 2.03(t, J=6Hz, 2H), 1.78(t, J=6Hz, 2H), 1.64(t, J=6Hz, 2H), 1.40(m, 6H); MS(FAB) m/z= 377.1(MH<sup>+</sup>-TFA).
- 25 30

## Example 21



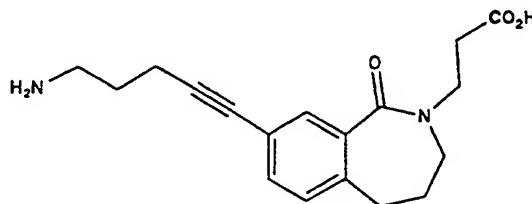
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**2-(2-carboxyethyl)-8-(8-guanadinooctyloxy)-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one trifluoroacetate.**

- a) A solution of 2-(2-carboxyethyl)-8-(8-aminooctyloxy)-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one trifluoroacetate (30 mg, 0.061 mmol) in saturated aqueous  $\text{KHCO}_3$  (2 mL) was treated with formamidine sulfonic acid (85 mg, 1.08 mmol) and the resulting mixture was stirred at room temperature for 20 minutes. The mixture was neutralized with acetic acid, and the residue was purified by HPLC (1/2' C-18 reverse-phase column) eluting with a gradient of 3:7 methanol (0.1% trifluoroacetate)/water (0.1% trifluoroacetate), time 0 to 10 min, to 7:3 methanol(0.1% trifluoroacetate)/water(0.1% trifluoroacetate), time 10 to 40 min, to yield 7.3 mg (25%) of 2-(2-carboxyethyl)-8-(8-guanadinooctyloxy)-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one trifluoroacetate as a light brown solid;  $^1\text{H}$  NMR( $\text{CD}_3\text{OD}$ )  $\delta$ 7.09(m, 2H), 6.96(dd,  $J=3, 9\text{Hz}$ , 1H), 3.98(t,  $J=6\text{Hz}$ , 2H), 3.83(t,  $J=6\text{Hz}$ , 2H), 3.33(m, 4H), 2.91(t,  $J=6\text{Hz}$ , 2H), 2.69(t,  $J=6\text{Hz}$ , 4H), 2.30(m, 2H), 1.75(m, 2H), 1.65(m, 2H), 1.40(m, 6H);  $^{13}\text{C}$  NMR( $\text{CD}_3\text{OD}$ )  $\delta$ 175.9(CO), 173.8(CO), 160.0, 137.0, 131.16, 130.79, 118.66, 114.81, 69.17, 45.33, 40.78, 34.32, 31.15, 30.27, 30.18, 30.15, 30.04, 28.59, 27.39, 27.06; MS(FAB)  $m/z$  = 419.2( $\text{MH}^+$ ).

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## Example 22



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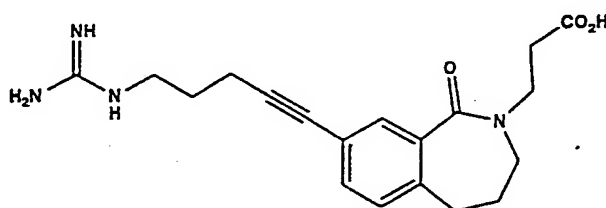
**2-(2-carboxyethyl)-8-(5-aminopentynyl)-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one trifluoroacetate.**

- a) A solution of 2-(2-carboxyethyl)-8-hydroxy-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one methyl ester (2.20 g, 0.00836 mol) and  $\text{NEt}_3$  (1.40 mL, 0.0100 mol) in  $\text{CH}_2\text{Cl}_2$  (20 mL) was cooled to  $0^\circ\text{C}$  (ice bath) and treated with trifluoromethanesulfonic anhydride (1.70 mL, 0.0100 mol), and the resulting mixture stirred to 30 minutes. The mixture was then partitioned between ethyl acetate and saturated aqueous  $\text{NaHCO}_3$ . The combined organic layers were

30

- dried ( $\text{MgSO}_4$ ), filtered, and concentrated *in vacuo* to an oil, used immediately without purification. A solution of the triflate (400 mg, 1.00 mmol) and 5-((*tert*-butyloxy)carbonyl)amino-2-pentyne (366 mg, 2.00 mmol) in  $\text{NEt}_3$  (10 mL) was treated with *bis*-triphenylphosphine palladium dichloride (20 mg) and  $\text{CuI}$  (10 mg), and the resulting mixture was de-gassed thoroughly with argon, and heated at reflux for 12 hours. The solution was then cooled, and volatiles removed *in vacuo*, and the residue was purified by chromatography over silica gel eluting with 1:1 ethyl acetate/hexanes to yield 400 mg (93%) of 2-(2-carboxyethyl)-8-(5-N-Boc-aminopentynyl)-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one methyl ester as a colorless oil;  $^1\text{H}$  NMR( $\text{CDCl}_3$ )  $\delta$  7.68(s, 1H), 7.37(d,  $J=6\text{Hz}$ , 1H), 7.50(d,  $J=6\text{Hz}$ , 1H), 4.7(brs, 1H), 3.84(t,  $J=6\text{Hz}$ , 2H), 3.71(s, 3H), 3.25(q,  $J=6\text{Hz}$ , 4H), 2.73(t,  $J=6\text{Hz}$ , 4H), 2.46(t,  $J=6\text{Hz}$ , 2H), 2.00(m, 2H), 1.78(m, 2H);  $^{13}\text{C}$  NMR( $\text{CDCl}_3$ )  $\delta$  171.38(CO), 170.49(CO), 155.92, 136.83, 136.06, 133.61, 131.68, 128.25, 122.55, 89.55, 80.43, 51.78, 47.18, 44.11, 39.75, 33.58, 29.92, 29.57, 28.72, 28.37, 16.88; MS(FAB)  $m/z$  = 429.1( $\text{MH}^+$ ).
- b) A solution of 2-(2-carboxyethyl)-8-(5-N-Boc-aminopentynyl)-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one methyl ester (66 mg, 0.154 mmol) in EtOH (1 mL) was treated with trifluoroacetic acid (0.5 mL), and the resulting mixture stirred for 18 hours. The volatiles were then removed, and the residue was dissolved in tetrahydrofuran (1 mL), and treated with 50% NaOH (0.5 mL), and the mixture was stirred for 20 minutes at room temperature. Trifluoroacetic acid was then added (3 mL), and the solution was concentrated *in vacuo*. The residue was purified by HPLC (1/2' C-18 reverse-phase column) eluting with a gradient of 3:7 methanol (0.1% trifluoroacetate)/water (0.1% trifluoroacetate), time 0 to 10 min, to 7:3 methanol(0.1% trifluoroacetate)/water(0.1% trifluoroacetate), time 10 to 40 min, to yield 46 mg (95%) of 2-(2-carboxyethyl)-8-(5-aminopentynyl)-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one trifluoroacetate as a light brown solid;  $^1\text{H}$  NMR( $\text{D}_2\text{O}$ )  $\delta$  7.32(s, 1H), 7.25(d,  $J=6\text{Hz}$ , 1H), 7.00(d,  $J=6\text{Hz}$ , 1H), 3.63(t,  $J=6\text{Hz}$ , 2H), 2.95(m, 4H), 2.58(t,  $J=6\text{Hz}$ , 2H), 2.50(t,  $J=6\text{Hz}$ , 2H), 2.38(t,  $J=6\text{Hz}$ , 2H), 1.80(m, 4H); MS(FAB)  $m/z$  = 315( $\text{MH}^+$ ).

## Example 23



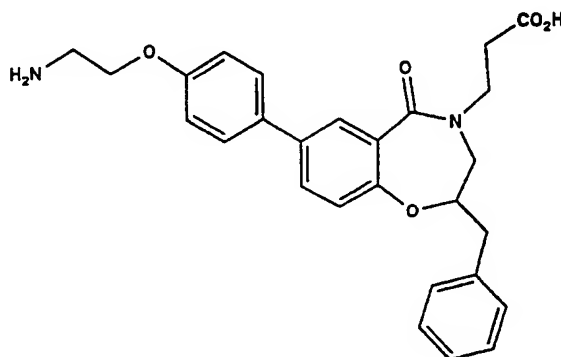
2-(2-carboxyethyl)-8-(5-guanidinopentynyl)-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one trifluoroacetate.

- a) A solution of 2-(2-carboxyethyl)-8-(5-aminopentynyl)-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one trifluoroacetate (20 mg, 0.063 mmol) in saturated aqueous  $\text{KHCO}_3$  (1 mL)

was treated with formamidine sulfonic acid (80 mg, 1.0 mmol), and the resulting mixture stirred for 2 hours. The mixture was then treated with acetic acid, and the volatiles were removed. The residue was purified by HPLC (1/2' C-18 reverse-phase column) eluting with a gradient of 3:7 methanol (0.1% trifluoroacetate)/water (0.1% trifluoroacetate), time 0 to 10 min, to 7:3 methanol(0.1% trifluoroacetate)/water(0.1% trifluoroacetate), time 10 to 40 min, to yield 4.5 mg (15%) of 2-(2-carboxyethyl)-8-(5-guanidinopentynyl)-2,3,4,5-tetrahydro-1H-2-benzazepin-1-one trifluoroacetate as a light brown solid;  $^1\text{H}$  NMR( $\text{D}_2\text{O}$ )  $\delta$ 7.40(s, 1H), 7.25(dd,  $J=3$ , 6Hz, 1H), 7.15(d,  $J=6$ Hz, 1H), 3.80(t,  $J=6$ Hz, 2H), 3.30(m, 4H), 3.15(t,  $J=6$ Hz, 2H), 2.00(t,  $J=6$ Hz, 2H), 1.80(m, 2H); MS(FAB)  $m/z=357.1(\text{MH}^+)$ .

10

## Example 24



15      2-(2-carboxyethyl)-8-(4-(2-aminoethoxy)phenyl)-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one trifluoroacetate.

a) A solution of 1-aminopropylene (10.6 grams, 0.186mol) in methanol was added to a methanolic solution of 3-phenylpropyleneoxide (5.07 grams, 0.037mol) and the mixture heated to reflux for 30 minutes. The mixture was allowed to warm to room temperature and concentrated *in vacuo* to yield 7.2 grams (99%) of N-allyl-3-phenyl-2-hydroxypropylamine.  $^1\text{H}$  NMR ( $\text{CDCl}_3$ ,  $\delta$ TMS) 7.34-7.18 (5H, m, Ar-H), 5.84 (1H, ddt,  $^3J_{\text{HH}}=17.1\text{Hz}$ ,  $^3J_{\text{HH}}=10.2\text{Hz}$ ,  $^3J_{\text{HH}}=6\text{Hz}$ ,  $\text{CH}_2\text{CH}$ ), 5.2-5.0 (2H, m,  $\text{CH}=\text{CH}_2$ ), 3.86 (1H, m,  $\text{CHOH}$ ), 3.19 (2H, m,  $\text{NCH}_2\text{CH}=\text{CH}_2$ ), 2.76-2.66 (3H, m,  $\text{NCH}_2\text{CH}(\text{OH})\text{CHHPh}$ ), 2.50 (1H, dd,  $^2J_{\text{HH}}=12\text{Hz}$ ,  $^3J_{\text{HH}}=9\text{Hz}$ ).  $^{13}\text{C}$  NMR ( $\text{CDCl}_3$ ) 138.4, 136.5, 129.3, 128.4, 126.3, 116.0, 70.6, 54.0, 51.9, 41.7.

25

Using the above procedure, but substituting the appropriate alkenyloxy or allyl amine for 3-phenylpropyleneoxide or 1-aminopropylene, respectively, there may be prepared, for example, the following compounds:

N-allyl-2-hydroxypropylamine, N-(3-but-1-ene)-3-phenyl-2-hydroxypropylamine, N-(3-but-1-ene)-2-hydroxypropylamine.

30

- b) To a magnetically stirred suspension of 5-iodosalicylic acid (Aldrich, 8.5 grams, 0.032 mol), was added oxalyl chloride (8.6 grams, 0.064 mol) and one drop of dimethylformamide. The mixture was stirred until effervescence had ceased and concentrated *in vacuo*. The resulting residue was diluted with diethyl ether and added to a solution of N-allyl-3-phenyl-2-hydroxypropyl-amine (6.4 grams, 0.034 mol) in diethyl ether over a period of 30 minutes. The reaction was stirred for 1 hour, filtered, and concentrated *in vacuo*. the resulting residue was dissolved in methylene chloride and washed with 1N HCl, sat. sodium bicarbonate, dried over sodium sulfate, filtered and concentrated in *vacuo* to yield 12.26 grams (87%) of N-allyl-N-(2-hydroxy-5-iodobenzoyl)-2-hydroxy-3-phenylpropylamine. The material was used without further purification.

Using the above procedure, but substituting the appropriate hydroxypropylamine for N-allyl-3-phenyl-2-hydroxypropylamine there may be prepared, for example, the following compounds:

- N-allyl-N-(2-hydroxy-5-iodobenzoyl)-2-hydroxypropylamine,  
N-(3-but-1-ene)-N-(2-hydroxy-5-iodobenzoyl)-2-hydroxy-3-phenylpropylamine,  
N-(3-but-1-ene)-N-(2-hydroxy-5-iodobenzoyl)-2-hydroxypropylamine.

- c) To a magnetically stirred solution of N-allyl-N-(2-hydroxy-5-iodobenzoyl)-2-hydroxy-3-phenylpropylamine (12.05 grams, 0.027 mol) in 100mL THF was added triphenylphosphine (7.24 grams, 0.027 mol), and diethyl azodicarboxylate (4.7 grams, 0.027 mol). The mixture was heated to reflux overnight, allowed to cool to room temperature, and concentrated *in vacuo*. the resulting residue was dissolved in ethyl acetate (10mL) and diluted with diethyl ether. The solids that form were filtered and the mother liquor was concentrated *in vacuo*. The yellow oil was purified by column chromatography (SiO<sub>2</sub>, using 15% diethyl ether/85% hexane to 1:1 diethyl ether/hexane as the eluting solvent gradient, R<sub>f</sub>(product, 1:1 Et<sub>2</sub>O/hexane)=0.64) to yield 4.31 grams (38%) of 2-allyl-4-benzyl-8-iodo-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one. <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 8.10 (1H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, C9-H), 7.69 (1H, dd, <sup>4</sup>J<sub>HH</sub>=2Hz, <sup>3</sup>J<sub>HH</sub>=9Hz, C7-H), 7.4-7.2 (5H, m, C<sub>6</sub>H<sub>5</sub>), 6.76 (1H, d, <sup>3</sup>J<sub>HH</sub>=9Hz, C6-H), 5.79 (1H, ddt, <sup>3</sup>J<sub>HH</sub>=17Hz, <sup>3</sup>J<sub>HH</sub>=11Hz, <sup>3</sup>J<sub>HH</sub>=6Hz, CH<sub>2</sub>CH=CH<sub>2</sub>), 5.2-5.0 (2H, m, CH=CH<sub>2</sub>), 5.71 (1H, m, CHOH), 4.31 (1H, ddt, <sup>2</sup>J<sub>HH</sub>=15Hz, <sup>3</sup>J<sub>HH</sub>=6Hz, <sup>4</sup>J<sub>HH</sub>=1Hz, NCHHCH=CH<sub>2</sub>), 4.00 (1H, ddt, <sup>2</sup>J<sub>HH</sub>=15Hz, <sup>3</sup>J<sub>HH</sub>=6Hz, <sup>4</sup>J<sub>HH</sub>=1Hz, NCHHCH=CH<sub>2</sub>), 3.34 (1H, dd, <sup>2</sup>J<sub>HH</sub>=16Hz, <sup>3</sup>J<sub>HH</sub>=5.1Hz, C3-H), 3.25 (1H, dd, <sup>2</sup>J<sub>HH</sub>=16Hz, <sup>3</sup>J<sub>HH</sub>=4.2Hz, C3-H), 3.09 (1H, dd, <sup>2</sup>J<sub>HH</sub>=14Hz, <sup>3</sup>J<sub>HH</sub>=7Hz, CHHPh), 2.73 (1H, dd, <sup>2</sup>J<sub>HH</sub>=14Hz, <sup>3</sup>J<sub>HH</sub>=7Hz, CHHPh). <sup>13</sup>C NMR (CDCl<sub>3</sub>) 166.9, 152.7, 141.4, 139.6, 136.5, 132.6, 129.8, 129.1, 128.8, 127.0, 124.6, 118.4, 86.7, 84.7, 50.4, 48.9, 38.9. IR (NaCl, cm<sup>-1</sup>) 1641, 1589, 1463, 1430, 1264, 1217, 699.

Using the above procedure, but substituting the appropriate hydroxypropylamine for N-allyl-N-(2-hydroxy-5-iodobenzoyl)-2-hydroxy-3-phenylpropylamine there may be prepared, for example, the following compounds:



2-allyl-4-methyl-8-iodo-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one,  
 2-(3-but-1-ene)-4-benzyl-8-iodo-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one,  
 2-(3-but-1-ene)-4-methyl-8-iodo-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one.

- 5 d) To a magnetically stirred solution of 2-allyl-4-benzyl-8-iodo-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one (0.9234 grams, 2.2mmol) in 4mL THF at room temperature under an atmosphere of nitrogen was added 6.0mL of a 0.5M solution of 9-borabicyclo[3.3.1]nonane in THF (3 mmol) and the mixture stirred for 1 hour. The reaction was quenched with 6mL of a solution of 2 grams of sodium hydroxide and 22mL of a 1:1 mixture of ethanol and water. To  
 10 the resulting solution was added 7mL of a 30% solution hydrogen peroxide in water. After 30 minutes the mixture was partitioned between diethyl ether and water. The organic layer was washed with 10% citric acid and brine, dried over sodium sulfate, concentrated *in vacuo*, and purified by column chromatography (SiO<sub>2</sub>, using 1:5 ethyl acetate/methylene chloride as the eluting solvent,  $R_f(\text{product})=0.22$ ) to yield 0.74 grams (77%) of 2-(3-hydroxypropyl)-4-benzyl-8-iodo-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one. <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 8.08 (1H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, C9-H), 7.71 (1H, dd, <sup>4</sup>J<sub>HH</sub>=2Hz, <sup>3</sup>J<sub>HH</sub>=8Hz, C7-H), 7.4-7.2 (5H, m, C<sub>6</sub>H<sub>5</sub>), 6.78 (1H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, C6-H), 4.64 (1H, m, ArOCH), 3.70 (2H, m, CH<sub>2</sub>OH), 3.58 (2H, bt, <sup>3</sup>J<sub>HH</sub>=5Hz, NCH<sub>2</sub>CH<sub>2</sub>), 3.39 (1H, dd, <sup>2</sup>J<sub>HH</sub>=15Hz, <sup>3</sup>J<sub>HH</sub>=8Hz, C3-H), 3.24 (1H, dd, <sup>2</sup>J<sub>HH</sub>=15Hz, <sup>3</sup>J<sub>HH</sub>=3Hz, C3-H), 3.1 (1H, dd, <sup>2</sup>J<sub>HH</sub>=14Hz, <sup>3</sup>J<sub>HH</sub>=7Hz, CHHPh), 2.76 (1H, dd, <sup>2</sup>J<sub>HH</sub>=14Hz, <sup>3</sup>J<sub>HH</sub>=7Hz, CHHPh), 1.68 (2H, m, CH<sub>2</sub>CH<sub>2</sub>OH). <sup>13</sup>C NMR (CDCl<sub>3</sub>) 168.3, 152.6, 141.6, 139.4, 136.2, 132.1, 131.9, 129.2, 129.0, 128.8, 128.4, 127.0, 124.7, 86.8, 85.7, 58.0, 49.9, 44.5, 38.8, 30.3. IR (NaCl, cm<sup>-1</sup>) 3422 (b), 1636, 1463, 1436, 1217, 1064.

Using the above procedure, but substituting the appropriate 5-oxa-2-benzazepin-1-one for 2-allyl-4-benzyl-8-iodo-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one there may  
 25 be prepared, for example, the following compounds:

2-(3-hydroxypropyl)-4-methyl-8-iodo-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one,  
 2-(4-hydroxybut-2-yl)-4-benzyl-8-iodo-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one,  
 2-(4-hydroxybut-2-yl)-4-methyl-8-iodo-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one.

- 30 e) To a magnetically stirred solution of 2-(3-hydroxypropyl)-4-benzyl-8-iodo-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one (0.75 grams, 1.7mmol) in 5mL dimethylformamide at room temperature was added a solution of pyridinium dichromate (2.58 grams, 6.86mmol) in 5mL dimethylformamide. The reaction mixture was stirred overnight, diluted with 75mL water, washed 3X75mL diethyl ether, made acidic to a pH of 2 with conc. HCl and extracted  
 35 3X75mL diethyl ether. The combined organics were extracted 3X40mL 1N sodium hydroxide. The combined aqueous layers were acidified to a pH of 2 with conc HCl and extracted 3X50mL diethyl ether. These ether layers were combined and dried over magnesium sulfate, filtered, and concentrated *in vacuo* to yield 0.557 grams (73%) of 2-(2-carboxyethyl)-4-benzyl-8-iodo-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one. <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 8.07 (1H, d,

<sup>4</sup>J<sub>HH</sub>=2Hz, C9-H), 7.69 (1H, dd, <sup>4</sup>J<sub>HH</sub>=2Hz, <sup>3</sup>J<sub>HH</sub>=8Hz, C7-H), 7.4-7.2 (5H, m, C<sub>6</sub>H<sub>5</sub>), 6.75 (1H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, C6-H), 4.75 (1H, m, ArOCH), 3.89 (1H, dt, <sup>2</sup>J<sub>HH</sub>=14Hz, <sup>3</sup>J<sub>HH</sub>=6Hz, NCHHCH<sub>2</sub>CO<sub>2</sub>), 3.69 (1H, dt, <sup>2</sup>J<sub>HH</sub>=14Hz, <sup>3</sup>J<sub>HH</sub>=6Hz, NCHHCH<sub>2</sub>CO<sub>2</sub>), 3.44 (2H, d, <sup>3</sup>J<sub>HH</sub>=3Hz, CH<sub>2</sub>Ph), 3.07 (1H, dd, <sup>2</sup>J<sub>HH</sub>=14Hz, <sup>3</sup>J<sub>HH</sub>=7Hz, C3-H), 2.8-2.7 (3H, m, C3-H, CH<sub>2</sub>CO<sub>2</sub>). <sup>13</sup>C NMR (CDCl<sub>3</sub>) 176.1, 167.6, 164.4, 152.8, 141.7, 139.5, 136.4, 129.4, 129.1, 128.9, 128.8, 127.0, 124.7, 86.7, 84.7, 51.4, 45.0, 38.8, 33.0.

Using the above procedure, but substituting the appropriate 5-oxa-2-benzazepin-1-one for 2-(3-hydroxypropyl)-4-benzyl-8-iodo-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one there may be prepared, for example, the following compounds:

- 10 2-(2-carboxyethyl)-4-methyl-8-iodo-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one,  
2-(3-carboxy-2-propyl)-4-benzyl-8-iodo-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one,  
2-(3-carboxy-2-propyl)-4-methyl-8-iodo-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one.

f) A solution of diazomethane in ether, prepared by allowing a slurry of 2.98 grams of N-nitroso-N-methylurea and 10mL ether to react with a solution of 20 grams of KOH in 20mL water, was carefully added *via* a pipet to solution of 2-(2-carboxyethyl)-8-iodo-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one (0.42 grams, 0.93mmol) in 20mL diethyl ether until yellow color persists. The mixture was allowed to stand overnight and concentrated *in vacuo*. The resulting residue was purified by column chromatography (SiO<sub>2</sub>, 1:9 ethyl acetate/hexane, R<sub>f</sub>(product)=0.2) to yield 0.275 grams (64%) of 2-(2-carboxyethyl)-8-iodo-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one methyl ester. <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 8.06 (1H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, C9-H), 7.68 (1H, dd, <sup>4</sup>J<sub>HH</sub>=2Hz, <sup>3</sup>J<sub>HH</sub>=8Hz, C7-H), 7.4-7.2 (5H, m, C<sub>6</sub>H<sub>5</sub>), 6.75 (1H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, C6-H), 4.76 (1H, m, ArOCH), 3.88 (1H, dt, <sup>2</sup>J<sub>HH</sub>=14Hz, <sup>3</sup>J<sub>HH</sub>=6Hz, NCHHCH<sub>2</sub>CO<sub>2</sub>), 3.69 (4H, m, NCHHCH<sub>2</sub>CO<sub>2</sub>CH<sub>3</sub>), 3.44 (2H, m, C3-H), 3.06 (1H, dd, <sup>2</sup>J<sub>HH</sub>=14Hz, <sup>3</sup>J<sub>HH</sub>=7Hz, C3-H), 2.76 (1H, dd, <sup>2</sup>J<sub>HH</sub>=14Hz, <sup>3</sup>J<sub>HH</sub>=7Hz, CHHPh), 2.69 (2H, m, CHHPh, CH<sub>2</sub>CO<sub>2</sub>). <sup>13</sup>C NMR (CDCl<sub>3</sub>) 172.4, 167.2, 164.4, 152.6, 141.4, 139.4, 136.5, 129.6, 129.1, 121.0, 128.8, 127.0, 124.6, 86.6, 84.7, 51.8, 51.3, 45.1, 38.7, 33.0.

Using the above procedure, but substituting the appropriate 5-oxa-2-benzazepin-1-one for 2-(2-carboxyethyl)-8-iodo-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one there may be prepared, for example, the following compounds:

- 30 2-(2-carboxyethyl)-4-methyl-8-iodo-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one methyl ester,  
2-(3-carboxy-2-propyl)-4-benzyl-8-iodo-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one methyl ester,  
35 2-(3-carboxy-2-propyl)-4-methyl-8-iodo-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one methyl ester.

g) To a magnetically stirred solution of 4-iodophenol (15 grams, 0.068mol) in 100mL THF was sequentially added triphenylphosphine (19.67 grams, 0.075mol), diethyl azodicarboxylate

(13.06 grams, 0.073mol), and 2-azidoethanol (5.93 grams, 0.068mol). The mixture was stirred overnight, concentrated in vacuo, diluted with 90mL ethyl acetate and diluted with hexane (400mL), and the resulting solids filtered away. The filtrate was concentrated in vacuo, diluted with diethyl ether, and extracted with 1N NaOH and water. Pentane was added to induce precipitation of a white solid that was again filtered away. The filtrate was concentrated *in vacuo*, and dissolved in methanol (60mL). Triethylamine (27 grams, 0.267mol) and propanedithiol (23 grams, 0.215mol) were added and the mixture stirred at room temperature overnight. The solution was diluted with 100mL water and partitioned between methylene chloride and 1N NaOH. The organic layer was extracted with 2X200mL 1N HCl and the aqueous layer washed with 100mL methylene chloride, made basic with 2N NaOH and extracted 3X75mL methylene chloride. The combined organics were dried over sodium carbonate, filtered and concentrated *in vacuo* to yield 12.97 grams of 2-(4-iodophenoxy)ethyl amine (72%). <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 7.56 (2H, d, <sup>3</sup>J<sub>HH</sub>=9Hz, Ar-H), 6.63 (2H, d, <sup>3</sup>J<sub>HH</sub>=9Hz, Ar-H), 4.03 (2H, t, <sup>3</sup>J<sub>HH</sub>=5Hz, CH<sub>2</sub>N<sub>3</sub>), 3.57 (2H, t, <sup>3</sup>J<sub>HH</sub>=5Hz, OCH<sub>2</sub>). <sup>13</sup>C NMR (CDCl<sub>3</sub>) 158.0, 138.2, 116.9, 83.4, 66.9, 49.9. IR (NaCl, cm<sup>-1</sup>) 2930, 2107 (s), 1589, 1483, 1284, 1237, 1058, 819.

h) 2-(4-iodophenoxy)ethylamine (10 grams, 0.038mol) dissolved in THF (50mL) was allowed to react with di-*t*-butyl dicarbonate (9.13 grams, 0.041mol) at room temperature for 1 hour. The reaction mixture was concentrated in vacuo to yield 17.7 grams (99%) of N-Boc-2-(4-iodophenoxy)ethylamine.

i) A slurry of 2-(2-carboxyethyl)-4-benzyl-8-iodo-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one methyl ester (0.155 grams, 0.334mmol), N-Boc-2-(4-iodophenoxy)ethylamine (0.246 grams, 0.664mmol), degassed triethylamine (2mL), bis(triphenylphosphine)palladium dichloride (11mgs, 10 mol %), and coprous iodide (6mgs, 20 mol %) was heated to reflux overnight. The reaction was concentrated *in vacuo*, triturating the resulting residue with diethyl ether (30mL). The ethereal solution was concentrated in vacuo and the residue purified by column chromatography (SiO<sub>2</sub>, using 1:1 ethyl acetate/hexane to ethyl acetate as the eluting solvent gradient, R<sub>f</sub>(product)=0.08) to yield 19.1mgs (10%) of 2-(2-carboxyethyl)-4-benzyl-8-(4-(2-N-Boc-amino)ethoxyphenyl)-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one methyl ester. <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 7.94 (1H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, C9-H), 7.59 (1H, dd, <sup>4</sup>J<sub>HH</sub>=2Hz, <sup>3</sup>J<sub>HH</sub>=8Hz, C7-H), 7.52 (2H, d, <sup>3</sup>J<sub>HH</sub>=9Hz, C8-*o*-Ar-H), 7.40-7.25 (5H, m, C<sub>6</sub>H<sub>5</sub>), 7.05 (1H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, C6-H), 6.96 (2H, d, <sup>3</sup>J<sub>HH</sub>=9Hz, C8-*m*-Ar-H), 5.03 (1H, bs, NH), 4.76 (1H, m, OCH), 4.06 (2H, t, <sup>3</sup>J<sub>HH</sub>=5Hz, OCH<sub>2</sub>CH<sub>2</sub>N), 3.92 (1H, dt, <sup>2</sup>J<sub>HH</sub>=14Hz, <sup>3</sup>J<sub>HH</sub>=7Hz, NCHHCH<sub>2</sub>CO<sub>2</sub>), 3.73 (1H, dt, <sup>2</sup>J<sub>HH</sub>=14Hz, <sup>3</sup>J<sub>HH</sub>=7Hz, NCHHCH<sub>2</sub>CO<sub>2</sub>), 3.67 (3H, s, CH<sub>3</sub>), 3.56 (2H, bq, <sup>3</sup>J<sub>HH</sub>=6Hz, CH<sub>2</sub>NH), 3.51 (1H, dd, <sup>2</sup>J<sub>HH</sub>=16Hz, <sup>3</sup>J<sub>HH</sub>=8Hz, C3-H), 3.43 (1H, dd, <sup>2</sup>J<sub>HH</sub>=16Hz, <sup>3</sup>J<sub>HH</sub>=5Hz, C3-H), 3.12 (1H, dd, <sup>2</sup>J<sub>HH</sub>=14Hz, <sup>3</sup>J<sub>HH</sub>=7Hz, CHHPh), 2.78 (1H, dd, <sup>2</sup>J<sub>HH</sub>=14Hz, <sup>3</sup>J<sub>HH</sub>=7Hz,

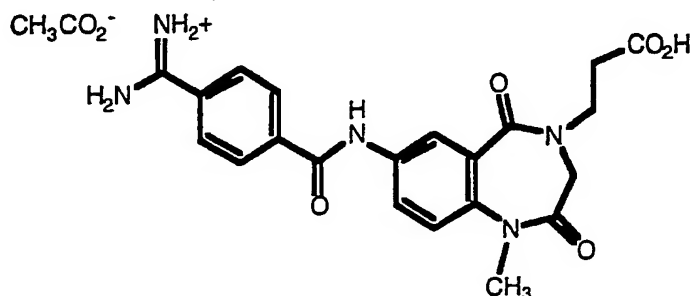
CHHPh), 2.72 (2H, m, CH<sub>2</sub>CO<sub>2</sub>), 1.48 (9H, s, Bu<sup>t</sup>), <sup>13</sup>C NMR (CDCl<sub>3</sub>) 172.6, 168.9, 136.8, 136.5, 130.8, 129.1, 128.8, 128.7, 128.3, 126.9, 123.0, 114.8, 84.6, 60.4, 51.8, 51.5, 45.0, 38.8, 33.2, 28.3.

- Using the above procedure, but substituting the appropriate 5-oxa-2-benzazepin-1-one for 2-(2-carboxyethyl)-8-iodo-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one there may be prepared, for example, the following compounds:
- 2-(2-carboxyethyl)-4-methyl-8-(4-(2-N-Boc-amino)ethoxyphenyl)-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one methyl ester,
- 2-(3-carboxy-2-propyl)-4-benzyl-8-(4-(2-N-Boc-amino)ethoxyphenyl)-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one methyl ester,
- 2-(3-carboxy-2-propyl)-4-methyl-8-(4-(2-N-Boc-amino)ethoxyphenyl)-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one methyl ester,
- 2-(2-carboxyethyl)-4-benzyl-8-[4-(4-N-Boc-piperazine)phenyl]-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one methyl ester,
- 2-(2-carboxyethyl)-4-methyl-8-[4-(4-N-Boc-piperazine)phenyl]-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one methyl ester,
- 2-(3-carboxy-2-propyl)-4-benzyl-8-[4-(4-N-Boc-piperazine)phenyl]-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one methyl ester,
- 2-(3-carboxy-2-propyl)-4-methyl-8-[4-(4-N-Boc-piperazine)phenyl]-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one methyl ester,
- 2-(2-carboxyethyl)-4-benzyl-8-[4-(4-N-Boc-aminomethyl)phenyl]-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one methyl ester,
- 2-(2-carboxyethyl)-4-methyl-8-[4-(4-N-Boc-aminomethyl)phenyl]-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one methyl ester,
- 2-(3-carboxy-2-propyl)-4-benzyl-8-[4-(4-N-Boc-aminomethyl)phenyl]-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one methyl ester,
- 2-(3-carboxy-2-propyl)-4-methyl-8-[4-(4-N-Boc-aminomethyl)phenyl]-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one methyl ester.
- j) To a solution of 2-(2-carboxyethyl)-4-benzyl-8-(4-(2-N-Boc-amino)ethoxyphenyl)-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one methyl ester (19.1mgs, 0.03mmol) in 0.5mL acetonitrile and 0.5mL water was added a concentrated solution of HCl in ethyl acetate (1mL) and the reaction allowed to stand for 15 minutes, concentrated *in vacuo*, dissolved in methanol (1mL) and allowed to react with 2N NaOH (1mL). The reaction was quenched with acetic acid (0.5mL) and concentrated *in vacuo*, dissolved in a minimum of water and purified by high-pressure column chromatography (HPLC, 1/2" reverse-phase C-18 column, 1:4 acetonitrile(0.1%TFA)/water (0.1%TFA) to 3:2 acetonitrile(0.1%TFA)/water (0.1%TFA) as the eluting solvent gradient) to yield 4.2mgs (30%) of 2-(2-carboxyethyl)-8-(4-(2-aminoethoxy)phenyl)-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one. <sup>1</sup>H NMR (CD<sub>3</sub>OD)

7.80 (1H, d,  $^4J_{HH}=2\text{Hz}$ , C9-H), 7.67 (1H, dd,  $^4J_{HH}=2\text{Hz}$ ,  $^3J_{HH}=8\text{Hz}$ , C7-H), 7.58 (2H, d,  $^3J_{HH}=9\text{Hz}$ , C8-*o*-Ar-H), 7.40-7.35 (5H, m, C<sub>6</sub>H<sub>5</sub>), 7.10 (2H, d,  $^3J_{HH}=9\text{Hz}$ , C8-*m*-Ar-H), 7.03 (1H, d,  $^3J_{HH}=8\text{Hz}$ , C6-H), 4.28 (2H, t,  $^3J_{HH}=4\text{Hz}$ , OCH<sub>2</sub>CH<sub>2</sub>), 3.81 (2H, m, NCH<sub>2</sub>CH<sub>2</sub>CO<sub>2</sub>), 3.61 (1H, dd,  $^2J_{HH}=14\text{Hz}$ ,  $^3J_{HH}=1\text{Hz}$ , C3-H), 3.44 (1H, dd,  $^2J_{HH}=14\text{Hz}$ ,  $^3J_{HH}=4\text{Hz}$ , C3-H), 3.41 (2H, t,  $^3J_{HH}=5\text{Hz}$ , CH<sub>2</sub>Ph), 2.87 (2H, t,  $^3J_{HH}=7\text{Hz}$ , CH<sub>2</sub>NH<sub>2</sub>), 2.50 (2H, t,  $^3J_{HH}=7\text{Hz}$ , CH<sub>2</sub>CO<sub>2</sub>). HRMS (FAB, MH<sup>+</sup>)  $m/z$  = 461.2109 (calcd for C<sub>27</sub>H<sub>29</sub>N<sub>2</sub>O<sub>5</sub>: 461.2076).

- 10           Using the above procedure, but substituting the appropriate 5-oxa-2-benzazepin-1-one for 2-(2-carboxyethyl)-4-benzyl-8-(4-(2-N-Boc-amino)ethoxyphenyl)-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one methyl ester there may be prepared, for example, the following compounds:
- 2-(2-carboxyethyl)-4-methyl-8-(4-(2-amino)ethoxyphenyl)-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one,
- 15   2-(3-carboxy-2-propyl)-4-benzyl-8-(4-(2-amino)ethoxyphenyl)-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one,
- 2-(3-carboxy-2-propyl)-4-methyl-8-(4-(2-amino)ethoxyphenyl)-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one,
- 20   2-(2-carboxyethyl)-4-benzyl-8-[4-(1-piperazine)phenyl]-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one,
- 2-(2-carboxyethyl)-4-methyl-8-[4-(1-piperazine)phenyl]-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one,
- 2-(3-carboxy-2-propyl)-4-benzyl-8-[4-(1-piperazine)phenyl]-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one,
- 25   2-(3-carboxy-2-propyl)-4-methyl-8-[4-(1-piperazine)phenyl]-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one,
- 2-(2-carboxyethyl)-4-benzyl-8-[4-(4-aminomethyl)phenyl]-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one,
- 30   2-(2-carboxyethyl)-4-methyl-8-[4-(4-aminomethyl)phenyl]-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one,
- 2-(3-carboxy-2-propyl)-4-benzyl-8-[4-(4-aminomethyl)phenyl]-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one,
- 2-(3-carboxy-2-propyl)-4-methyl-8-[4-(4-aminomethyl)phenyl]-2,3,4,5-tetrahydro-1H-5-oxa-2-benzazepin-1-one.
- 35

## Example 25



5                    **1-Methyl-4-(2-carboxyethyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione acetate**

a) A solution of N-methylisatoic anhydride (64.0 g, 0.325 mol) and b-alanine ethyl ester hydrochloride (50.0 g, 0.325 mol) in N,N-dimethylformamide (325 mL) was treated with  
 10 N,N-dimethylaminopyridine (2.0 g), the resulting mixture cooled to 0°C (ice bath), and triethylamine (46.0 mL, 0.325 mol) was added in one batch. Vigorous CO<sub>2</sub> evolution was noted. The mixture was stirred for 18 h at room temperature, then diluted with 400 mL ethyl acetate. The solution was washed exhaustively with water and brine to remove the N,N-dimethylformamide, then dried over MgSO<sub>4</sub>, filtered and concentrated in vacuo. The residue  
 15 was adsorbed on silica gel, and elution with a gradient of hexane to 1:1 ethyl acetate-hexane yielded 80 g (98%) of an oil. <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 7.28 (3H, m), 6.75 (1H, s, NHCH<sub>3</sub>), 6.62 (1H, d, J=9Hz, ArH), 6.55 (1H, t, J=9Hz, ArH), 4.14 (q, J=9Hz, 2H, CH<sub>2</sub>O), 3.63 (q, J=6Hz, NCH<sub>2</sub>CH<sub>2</sub>), 2.83 (s, 3H, NCH<sub>3</sub>), 2.59 (2H, t, J=6Hz, CH<sub>2</sub>CO<sub>2</sub>Et), 1.24 (3H, t, J=9Hz, CH<sub>2</sub>CH<sub>3</sub>); <sup>13</sup>C NMR (CDCl<sub>3</sub>, dTMS) 172.88, 169.89, 150.71, 132.96, 127.42, 115.01, 114.60,  
 20 111.16, 60.87, 35.11, 34.19, 29.74, 14.28.

b) A solution of N-(N-methyl-2-aminobenzoyl)-b-alanine ethyl ester (80.0 g, 0.319 mol) in CH<sub>2</sub>Cl<sub>2</sub> (400 mL) and H<sub>2</sub>O (200 mL) was cooled to 0°C (ice bath) and treated rapidly  
 25 dropwise with bromoacetyl bromide (42 mL, 0.478 mol). After the addition was complete, the phases were separated, and the aqueous phase was washed twice with CH<sub>2</sub>Cl<sub>2</sub>, and the combined organic phases were dried (MgSO<sub>4</sub>), filtered and concentrated *in vacuo* to yield 105 g (89%) of an oil. This material was dissolved in N,N-dimethylformamide (200 mL), and with rapid stirring, was treated in one portion with Cs<sub>2</sub>CO<sub>3</sub> (93.0 g, 0.283 mol). The mixture was stirred for 1/2 h at room temperature, then diluted with ethyl acetate and washed  
 30 exhaustively with brine and water to remove N,N-dimethylformamide. The solution was then dried (MgSO<sub>4</sub>), filtered and concentrated in vacuo. The residue was dissolved in minimal ethyl acetate and diluted with hexane until cloudy. A small amount of ethyl acetate was then added, and the mixture stored at -20°C. The crystals that had formed were

collected and washed with hexane, then dried *in vacuo* to yield 43 g (52%) of colorless crystals, mp. 119-121°C. <sup>1</sup>H NMR(CDCl<sub>3</sub>, dTMS) 7.83 (1H, dd, J=1.5, 8Hz, ArH), 7.51 (1H, td, J=1.5, 6Hz, ArH), 7.28 (1H, t, J=6Hz, ArH), 7.18 (1H, d, J=9Hz, ArH), 4.13(2H, q, J=9Hz, CH<sub>2</sub>O), 4.03 (1H, B part, ABq, J=15Hz, CH<sub>2</sub>CO), 3.92 (2H, t, J=6Hz, NCH<sub>2</sub>), 3.84 (1H, A part, ABq, J=15Hz, CH<sub>2</sub>CO), 3.37 (s, 3H, NCH<sub>3</sub>), 2.71 (2H, m, complex, CH<sub>2</sub>CO<sub>2</sub>Et), 1.24 (3H, t, J=6Hz, CH<sub>2</sub>CH<sub>3</sub>), 171.43, 169.16, 167.26, 141.08, 132.25, 130.92, 128.74, 125.75, 121.07, 60.89, 52.31, 45.18, 35.03, 32.97, 14.29; MS(FAB) *m/z* 291(MH<sup>+</sup>).

c) Fuming nitric acid (200 mL) was cooled to 0°C and treated portionwise with 1-methyl-4-(2-carboxyethyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester (43.0 g, 0.148 mol), and the resulting dark yellow mixture was stirred for 24 h at room temperature. The solution was then poured over ice onto solid NaHCO<sub>3</sub> and extracted exhaustively with ethyl acetate. The combined organic layers were dried (MgSO<sub>4</sub>), filtered and concentrated *in vacuo* to yield 36 g (73% ) of an orange oil; <sup>1</sup>H NMR(CDCl<sub>3</sub>, dTMS) 8.70 (1H, d, J=3Hz, ArH), 8.34 (1H, dd, J=3, 9Hz, ArH), 7.38 (1H, d, J=9Hz, ArH), 4.14 (2H, q, J=6Hz, CH<sub>2</sub>O), 4.04 (2H, d, J=6Hz, CH<sub>2</sub>CO), 3.93 (2H, t, J=6Hz, NCH<sub>2</sub>), 3.44 (3H, s, NCH<sub>3</sub>), 2.74 (2H, m, complex, CH<sub>2</sub>CO<sub>2</sub>Et), 1.26 (3H, t, J=6Hz, CH<sub>2</sub>CH<sub>3</sub>); <sup>13</sup>C NMR(CDCl<sub>3</sub>) d194.51, 191.54, 188.52, 168.96, 167.69, 152.58, 150.31, 150.05, 145.12, 84.22, 75.21, 68.61, 58.39, 55.97, 37.44; MS(FAB) *m/z* 336.1(MH<sup>+</sup>).

Using the same above method, but substituting the appropriate 1-alkyl-4-(2-carboxyethyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester for 1-methyl-4-(2-carboxyethyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester there may be prepared, for example, the following compounds:

- 1-ethyl-4-(2-carboxyethyl)-7-nitro-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 25 1-isopropyl-4-(2-carboxyethyl)-7-nitro-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-isobutyl-4-(2-carboxyethyl)-7-nitro-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-cyclohexyl-4-(2-carboxyethyl)-7-nitro-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester.
- 30

d) A solution of 1-methyl-4-(2-carboxyethyl)-7-nitro-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester (36.0 g, .107 mol) in ethyl acetate (200 mL) and absolute ethanol (150 mL) was hydrogenated over Pd on carbon (4 g) at 40 PSI for 24 h. The black suspension was then filtered and concentrated *in vacuo* to give 32.3 g (100%) of a colorless foam. This foam was dissolved in CH<sub>2</sub>Cl<sub>2</sub>, cooled to 0°C (ice bath) and treated with NEt<sub>3</sub> (17.5 mL, 0.126 mol) followed by 4-cyanobenzoyl chloride (20.86 g, 0.126 mol), and the resulting mixture was stirred for 24 h at room temperature. The mixture was then washed with water and brine, dried over MgSO<sub>4</sub>, filtered and concentrated *in vacuo* to yield a light brown solid.

35

Recrystallization of the solid from  $\text{CH}_2\text{Cl}_2$ /hexanes yielded 33 g (76%) of a white granular solid: mp. 158-160°C;  $^1\text{H}$  NMR( $\text{CDCl}_3$ , dTMS) 8.76 (1H, s, NH), 8.00 (1H, dd,  $J=3$ , 9Hz, ArH), 7.72 (1H, d,  $J=9$ Hz, ArH), 7.58 (1H, d,  $J=3$ Hz, ArH), 7.47 (1H, d,  $J=9$ Hz, ArH), 6.94 (1H, d,  $J=9$ Hz, ArH), 3.83 (2H, q,  $J=6$ Hz,  $\text{OCH}_2\text{CH}_3$ ), 3.77 (1H, B part, ABq,  $J=15$ Hz,  $\text{CH}_2\text{CO}$ ), 3.57 (1H, A part, ABq,  $J=15$ Hz,  $\text{CH}_2\text{CO}$ ), 3.53 (2H, t,  $J=6$ Hz,  $\text{NCH}_2$ ), 3.09 (3H, s,  $\text{NCH}_3$ ), 2.34 (2H, m, complex,  $\text{CH}_2\text{CO}_2\text{Et}$ ), 0.95 (3H, t,  $J=6$ Hz,  $\text{OCH}_2\text{CH}_3$ );  $^{13}\text{C}$  NMR( $\text{CDCl}_3$ ) d171.04, 168.57, 166.90, 164.31, 138.24, 137.34, 135.51, 132.47, 128.60, 128.05, 124.64, 121.85, 117.83, 115.51, 60.94, 52.09, 45.01, 34.92, 32.60, 14.13; MS(FAB)  $m/z$  435( $\text{MH}^+$ ).

- Using the same above procedure, but substituting the appropriate cyano acid chloride or amino protected amino acid chloride for 4-cyanobenzoyl chloride there may be prepared, for example, the following compounds:
- 1-methyl-4-(2-carboxyethyl)-7-(3-cyano)benzamido-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-methyl-4-(2-carboxyethyl)-7-(4-cyano)butanamido-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-methyl-4-(2-carboxyethyl)-7-(5-cyano)pentamido-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-methyl-4-(2-carboxyethyl)-7-4-(N-Boc)-aminobutanamido-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-methyl-4-(2-carboxyethyl)-7-5-(N-Boc)-aminopentamido-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester. Furthermore, using the same procedure but substituting the appropriate 1-alkyl-4-(2-carboxyethyl)-7-nitro-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester for 1-methyl-4-(2-carboxyethyl)-7-nitro-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester there may be prepared, for example, the following compounds:
- 1-ethyl-4-(2-carboxyethyl)-(4-cyano)benzamido-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-isopropyl-4-(2-carboxyethyl)-(4-cyano)benzamido-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-isobutyl-4-(2-carboxyethyl)-(4-cyano)benzamido-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-cyclohexyl-4-(2-carboxyethyl)-(4-cyano)benzamido-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester.

- e) A solution of 1-methyl-4-(2-carboxyethyl)-7-(4-cyano)benzamido-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester (10.0 g, 0.023 mol) in  $\text{NEt}_3$  (50 mL) and pyridine (70 mL) was saturated with  $\text{H}_2\text{S}$ , then heated at 70°C for 24 h. The solvents were removed in vacuo, and the residue remaining was suspended in  $\text{CH}_2\text{Cl}_2$  (200 mL) and treated with iodomethane (10 mL). The resulting solution was heated at reflux for 24 h. The volatiles were then



removed in *vacuo*, and the resulting residue was dissolved in EtOH (100 mL) and treated with  $\text{NH}_4\text{OAc}$  (3.2 g), and the resulting suspension heated to  $50^\circ\text{C}$  with stirring for 24 h. The mixture was concentrated in *vacuo*, and the product was dissolved in 0.5% aqueous  $\text{HOAc}$ , and isolated by HPLC to yield a white solid;  $^1\text{H}$  NMR ( $\text{D}_2\text{O}$ , TFA salt)  $\delta$  7.72 (2H, d,  $J=12\text{Hz}$ , ArH), 7.65 (2H, d,  $J=12\text{Hz}$ , ArH), 7.63 (1H, d,  $J=3\text{Hz}$ , ArH), 7.53 (1H, dd,  $J=3, 9\text{Hz}$ , ArH), 7.12 (1H, d,  $J=12\text{Hz}$ , ArH), 4.40 (2H, B part, ABq,  $J=15\text{Hz}$ ,  $\text{C}_3\text{-H}$  and m,  $\text{CHHCH}_2\text{CO}_2$ ), 3.91 (2H, q,  $J=6\text{Hz}$ ,  $\text{OCH}_2$ ), 3.68 (1H, A part, ABq,  $\text{C}_3\text{-H}$ ), 3.48 (1H, m,  $\text{CHHCH}_2\text{CO}_2$ ), 3.18 (3H, s,  $\text{NCH}_3$ ), 2.54 (2H, m,  $\text{CH}_2\text{CO}_2$ ), 0.98 (3H, t,  $J=6\text{Hz}$ ,  $\text{OCH}_2\text{CH}_3$ ); MS (FAB)  $m/z$  452 ( $\text{MH}^+$ ).

- Using the above procedure, but substituting the appropriate 1-alkyl-4-(2-carboxyethyl)-7-(4-cyano)alkyl/benzamido-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester for 1-methyl-4-(2-carboxyethyl)-7-(4-cyano)benzamido-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester, there may be prepared, for example, the following compounds.
- 10 1-methyl-4-(2-carboxyethyl)-7-(3-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
  - 15 1-methyl-4-(2-carboxyethyl)-7-(4-amidino)butanamido-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
  - 1-methyl-4-(2-carboxyethyl)-7-(5-amidino)pentamido-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
  - 20 1-methyl-4-(2-carboxyethyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
  - 1-ethyl-4-(2-carboxyethyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
  - 25 1-isopropyl-4-(2-carboxyethyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
  - 1-isobutyl-4-(2-carboxyethyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
  - 1-cyclohexyl-4-(2-carboxyethyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester.

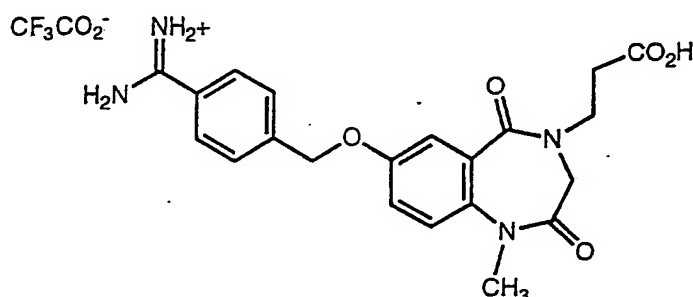
- f) A solution of 1-methyl-4-(2-carboxyethyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester acetate (0.2 g, 0.456 mmol) in THF (20 mL) and 50% aqueous  $\text{NaOH}$  (3 mL) was stirred for 24 h at room temperature. Acetic acid was then added, and the solution was concentrated in *vacuo*. The residue remaining was purified by HPLC to yield 122 mg (57%) of a white foam (isolated by lyophilization);  $^1\text{H}$  NMR ( $\text{D}_2\text{O}$ )  $\delta$  7.73 (2H, d,  $J=12\text{Hz}$ , ArH), 7.63 (2H, d,  $J=6\text{Hz}$ , ArH), 7.62 (1H, m, ArH), 7.54 (1H, dd,  $J=3, 9\text{Hz}$ , ArH), 7.12 (1H, d,  $J=12\text{Hz}$ , ArH), 3.87 (1H, B part, ABq,  $J=15\text{Hz}$ ,  $\text{CH}_2\text{CO}$ ) overlapping 3.84 (1H, m, complex), 3.20 (1H, A part, ABq,  $J=15\text{Hz}$ ,  $\text{CH}_2\text{CO}$ ), 3.48 (1H, m), 3.03 (3H, s,  $\text{NCH}_3$ ), 2.50 (2H, m, complex,  $\text{CH}_2\text{CO}$ ); MS (FAB)  $m/z$  424.3 ( $\text{MH}^+$ ); High Resolution MS (FAB) calculated for  $\text{C}_{21}\text{H}_{22}\text{N}_5\text{O}_5$  424.1620, found 424.1647.

Using the above method, but substituting the appropriate 1-alkyl-4-(2-carboxyethyl)-7-(4-amidino)alkyl/benzamido-3,4-dihydro-1H-1,4-benzodiazepine-2,5-

dione ethyl ester for 1-methyl-4-(2-carboxyethyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester, there may be prepared, for example, the following compounds:

- 1-methyl-4-(2-carboxyethyl)-7-(3-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione acetate,
- 1-methyl-4-(2-carboxyethyl)-7-(4-amidino)butanamido-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione acetate,
- 1-methyl-4-(2-carboxyethyl)-7-(5-amidino)pentamido-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione acetate,
- 1-ethyl-4-(2-carboxyethyl)-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione acetate,
- 1-isopropyl-4-(2-carboxyethyl)-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione acetate,
- 1-isobutyl-4-(2-carboxyethyl)-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione acetate,
- 1-cyclohexyl-4-(2-carboxyethyl)-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione acetate.

Example 26



*1-methyl-4-(2-carboxyethyl)-7-(4-amidinobenzoyloxy)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate.*

25

- a) To a stirred solution of 643 mg (2.11 mmol) of 1-methyl-4-(2-carboxyethyl)-7-amino-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester (prepared from 1-methyl-4-(2-carboxyethyl)-7-nitro-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester as a colorless foam as described in part (d) of example 26) in 60 mL of 0.1 N sulfuric acid at 0 °C was added a solution of 153 mg (2.21 mmol) of sodium nitrite in 5 ml of water. After 10 min., a solution of 49.0 g (211 mmol) of copper (II) nitrate hemipentahydrate in 90 mL of water was added followed by 287 mg (2.00 mmol) of copper (II) oxide. The mixture was warmed to ambient temperature and extracted with three portions of methylene chloride. The combined

extracts were washed with water, brine, dried over anhydrous magnesium sulfate, and concentrated *in vacuo*. Chromatography (silica gel 60, gradient elution from 1:2 ethyl acetate/hexane to ethyl acetate,  $R_f = 0.2$  in 2:1 ethyl acetate/hexane on silica) gave 353 mg (55%) of 1-methyl-4-(2-carboxyethyl)-7-hydroxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester as a colorless stable foam.  $^1\text{H}$  NMR (300 MHz,  $\text{CDCl}_3$ )  $\delta$  8.60 (1H, bs, Ar-OH), 7.52 (1H, bs, Ar-H), 7.1-7.0 (2H, m, Ar-H), 4.12 (2H, q,  $J=7.3$  Hz,  $\text{OCH}_2\text{CH}_3$ ), 4.07 (1H, d,  $J=14.7$  Hz,  $\text{NCHHCON}$ ), 3.90 (2H, m,  $\text{NCH}_2\text{CH}_2$ ), 3.86 (1H, d,  $J=14.7$  Hz,  $\text{NCHHCON}$ ), 3.33 (3H, s,  $\text{NCH}_3$ ), 2.70 (2H, m,  $\text{NCH}_2\text{CH}_2$ ), 1.23 (3H, t,  $J=7.3$  Hz,  $\text{OCH}_2\text{CH}_3$ ); Exact mass (FAB,  $\text{M}+\text{H}^+$ ) calcd for  $\text{C}_{15}\text{H}_{19}\text{N}_2\text{O}_5$ : 307.1294; Found: 307.1289.

- 10 Using the above procedure, but substituting the appropriate 1-alkyl-4-(2-carboxyethyl)-7-nitro-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester for 1-methyl-4-(2-carboxyethyl)-7-nitro-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester, there can be prepared, for example, the following compounds:
- 1-ethyl-4-(2-carboxyethyl)-7-hydroxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 15 1-isopropyl-4-(2-carboxyethyl)-7-hydroxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 1-isobutyl-4-(2-carboxyethyl)-7-hydroxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 20 1-cyclohexyl-4-(2-carboxyethyl)-7-hydroxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester.

- b) To a stirred solution of 121 mg (0.395 mmol) of 1-methyl-4-(2-carboxyethyl)-7-hydroxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester and 93.0 mg (0.474 mmol) of  $\alpha$ -bromo-p-tolynitrile in 2 ml of dry dimethylformamide was added 60.0 mg (0.434 mmol) of potassium carbonate. The mixture was stirred overnight at ambient temperature then partitioned between ethyl acetate and water. The organic phase was separated and washed with water, brine, dried over anhydrous magnesium sulfate, and concentrated *in vacuo*. Chromatography (silica gel 60, gradient elution from 2:1 ethyl acetate/hexane to ethyl acetate,  $R_f = 0.35$  in 2:1 ethyl acetate/hexane on silica) gave 125 mg (75%) of 1-methyl-4-(2-carboxyethyl)-7-(4-cyanobenzyloxy)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester as a colorless foam.  $^1\text{H}$  NMR (300 MHz,  $\text{CDCl}_3$ )  $\delta$  7.68 (2H, d,  $J=8.8$  Hz, NC-Ar-H<sub>2</sub>), 7.54 (2H, d,  $J=8.8$  Hz, NC-Ar-H<sub>2</sub>), 7.52 (1H, bs, O-Ar-H), 7.13 (2H, bs, O-Ar-H), 5.16 (2H, ABq,  $J=12.7$  Hz,  $\text{ArCH}_2\text{O}$ ), 4.13 (2H, q,  $J=7.3$  Hz,  $\text{OCH}_2\text{CH}_3$ ), 4.03 (1H, d,  $J=14.7$  Hz,  $\text{NCHHCON}$ ), 3.92 (2H, m,  $\text{NCH}_2\text{CH}_2$ ), 3.84 (1H, d,  $J=14.7$  Hz,  $\text{NCHHCON}$ ), 3.34 (3H, s,  $\text{NCH}_3$ ), 2.70 (2H, m,  $\text{NCH}_2\text{CH}_2$ ), 1.24 (3H, t,  $J=7.3$  Hz,  $\text{OCH}_2\text{CH}_3$ ); Exact mass (FAB,  $\text{M}+\text{H}^+$ ) calcd for  $\text{C}_{23}\text{H}_{24}\text{N}_3\text{O}_5$ : 422.1716; Found: 422.1744.

Using the above procedure, but substituting the appropriate 1-alkyl-4-(2-carboxyethyl)-7-hydroxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester for 1-methyl-4-(2-carboxyethyl)-7-hydroxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester there may be prepared, for example, the following compounds:

- 5 1-ethyl-4-(2-carboxyethyl)-7-(4-cyanobenzyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,  
1-isopropyl-4-(2-carboxyethyl)-7-(4-cyanobenzyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,  
1-isobutyl-4-(2-carboxyethyl)-7-(4-cyanobenzyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 10 1-cyclohexyl-4-(2-carboxyethyl)-7-(4-cyanobenzyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,  
1-ethyl-4-(2-carboxyethyl)-7-(3-cyanobenzyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 15 1-isopropyl-4-(2-carboxyethyl)-7-(3-cyanobenzyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,  
1-isobutyl-4-(2-carboxyethyl)-7-(3-cyanobenzyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,  
1-cyclohexyl-4-(2-carboxyethyl)-7-(3-cyanobenzyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 20 1-ethyl-4-(2-carboxyethyl)-7-(4-cyanobutyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,  
1-isopropyl-4-(2-carboxyethyl)-7-(4-cyanobutyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,
- 25 1-isobutyl-4-(2-carboxyethyl)-7-(4-cyanobutyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,  
1-cyclohexyl-4-(2-carboxyethyl)-7-(4-cyanobutyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester.

- 30 c) Hydrogen sulfide gas was bubbled into a solution of 83.6 mg (0.198 mmol) of 1-methyl-4-(2-carboxyethyl)-7-(4-cyanobenzyloxy)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester in 3 mL of 1:1 pyridine/triethylamine for 15 min and the mixture was warmed in a 50 °C oil bath for 4 h. After evaporation of the solvent, the residue was taken up in 6 mL of 5:1 methylene chloride/methyl iodide and warmed at 50 °C for 1 h. The solution was again
- 35 concentrated and the residue redissolved in 3 mL of methanol containing 0.5 g of ammonium acetate. The mixture was heated at 50 °C for 12 h, concentrated to a volume of 1 mL, and purified by reverse phase HPLC (Rainin Microsorb 1/2" C-18). The product was eluted with a solvent gradient of 0:100 acetonitrile (0.1% trifluoroacetic acid) /water (0.1% trifluoroacetic acid), time 0 to 10 min, to 50:50 acetonitrile (0.1% trifluoroacetic acid)/water (0.1%

trifluoroacetic acid), time 10 to 40 min, flow=12 ml/min,  $R_t$ =35.2 min, uv detection at 254 nm.

Concentration of the clean fractions gave 70.0 mg (65%) of 1-methyl-4-(2-carboxyethyl)-7-(4-amidinobenzoyloxy)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester

trifluoroacetate as a colorless amorphous solid.  $^1\text{H}$  NMR (300 MHz, 1:1  $\text{D}_2\text{O}/\text{d}^6\text{-acetone}$ ) d

- 5 7.85 (2H, d,  $J=7.5$  Hz,  $\text{HNC-Ar-H}_2$ ), 7.68 (2H, d,  $J=7.5$  Hz,  $\text{HNC-Ar-H}_2$ ), 7.40-7.25 (3H, m, O-Ar-H<sub>3</sub>), 5.24 (2H, s,  $\text{ArCH}_2\text{O}$ ), 4.07 (1H, d,  $J=15.1$  Hz,  $\text{NCHHCON}$ ), 4.04 (1H, m,  $\text{NCHHCH}_2$ ), 4.02 (2H, q,  $J=7.3$  Hz,  $\text{OCH}_2\text{CH}_3$ ), 3.82 (1H, d,  $J=15.1$  Hz,  $\text{NCHHCON}$ ), 3.71 (1H, ddd,  $J=14.2, 5.9, 5.9$  Hz,  $\text{NCHHCH}_2$ ), 3.29 (3H, s,  $\text{NCH}_3$ ), 2.65 (2H, m,  $\text{NCH}_2\text{CH}_2$ ), 1.11 (3H, t,  $J=7.3$  Hz,  $\text{OCH}_2\text{CH}_3$ ); Exact mass (FAB,  $\text{M}+\text{H}^+$ ) calcd for  $\text{C}_{23}\text{H}_{27}\text{N}_4\text{O}_5$ : 439.1981;
- 10 Found: 439.1977.

Using the above procedure, but substituting the appropriate 1-alkyl-4-(2-carboxyethyl)-7-(cyanobenzyl/alkyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester for 1-methyl-4-(2-carboxyethyl)-7-(4-cyanobenzyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester there may be prepared, for example, the following

- 15 compounds:

1-ethyl-4-(2-carboxyethyl)-7-(4-amidinobenzyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,

1-isopropyl-4-(2-carboxyethyl)-7-(4-amidinobenzyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,

- 20 1-isobutyl-4-(2-carboxyethyl)-7-(4-amidinobenzyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,

1-cyclohexyl-4-(2-carboxyethyl)-7-(4-amidinobenzyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,

- 25 1-ethyl-4-(2-carboxyethyl)-7-(3-amidinobenzyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,

1-isopropyl-4-(2-carboxyethyl)-7-(3-amidinobenzyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,

- 30 1-isobutyl-4-(2-carboxyethyl)-7-(3-amidinobenzyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,

1-cyclohexyl-4-(2-carboxyethyl)-7-(3-amidinobenzyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,

- 35 1-ethyl-4-(2-carboxyethyl)-7-(4-amidinobutyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,

1-isopropyl-4-(2-carboxyethyl)-7-(4-amidinobutyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,

- 1-isobutyl-4-(2-carboxyethyl)-7-(4-amidinobutyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,

1-cyclohexyl-4-(2-carboxyethyl)-7-(4-amidinobutyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester.

- d) To a stirred, 0 °C solution of 37.3 mg (0.0690 mmol) of 1-methyl-4-(2-carboxyethyl)-7-(4-amidinobenzoyloxy)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester trifluoroacetate in 1 mL of tetrahydrofuran was added 0.28 mL of 1 N aqueous sodium hydroxide and the mixture warmed to ambient temperature over 1 h. The reaction was acidified with 50 mL of acetic acid, concentrated in vacuo, and purified by reverse phase HPLC (Rainin Microsorb 1/2" C-18). The product was eluted with a solvent gradient of 0:100 acetonitrile (0.1% trifluoroacetic acid) /water (0.1% trifluoroacetic acid), time 0 to 10 min, to 50:50 acetonitrile (0.1% trifluoroacetic acid)/water (0.1% trifluoroacetic acid), time 10 to 40 min, flow=12 ml/min,  $R_t$ =32.4 min, uv detection at 254 nm. Concentration of the clean fractions gave 26.6 mg (75%) of 1-methyl-4-(2-carboxyethyl)-7-(4-amidinobenzoyloxy)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate as a colorless amorphous solid.  $^1\text{H}$  NMR (300 MHz, 1:1  $\text{D}_2\text{O}/\text{d}^6$ -acetone)  $\delta$  7.75 (2H, d,  $J$ =8.3 Hz, HNC-Ar-H<sub>2</sub>), 7.60 (2H, d,  $J$ =8.3 Hz, HNC-Ar-H<sub>2</sub>), 7.32-7.18 (3H, m, O-Ar-H<sub>3</sub>), 5.16 (2H, s, ArCH<sub>2</sub>O), 4.01 (1H, d,  $J$ =15.1 Hz, NCHHCON), 4.04 (1H, ddd,  $J$ =14.2, 7.3, 7.3 Hz, NCHHCH<sub>2</sub>), 3.77 (1H, d,  $J$ =15.1 Hz, NCHHCON), 3.67 (1H, ddd,  $J$ =14.2, 5.8, 5.8 Hz, NCHHCH<sub>2</sub>), 3.20 (3H, s, NCH<sub>3</sub>), 2.59 (2H, m, NCH<sub>2</sub>CH<sub>2</sub>); Exact mass (FAB,  $M+H^+$ ) calcd for  $\text{C}_{21}\text{H}_{23}\text{N}_4\text{O}_5$ : 411.1668; Found: 411.1701.

Using the above procedure, but substituting the appropriate 1-alkyl-4-(2-carboxyethyl)-7-(amidinobenzyl/alkyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester for 1-methyl-4-(2-carboxyethyl)-7-(4-amidinobenzyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester there may be prepared, for example, the following compounds:

- 1-ethyl-4-(2-carboxyethyl)-7-(4-amidinobenzyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-isopropyl-4-(2-carboxyethyl)-7-(4-amidinobenzyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-isobutyl-4-(2-carboxyethyl)-7-(4-amidinobenzyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-cyclohexyl-4-(2-carboxyethyl)-7-(4-amidinobenzyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-ethyl-4-(2-carboxyethyl)-7-(3-amidinobenzyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-isopropyl-4-(2-carboxyethyl)-7-(3-amidinobenzyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-isobutyl-4-(2-carboxyethyl)-7-(3-amidinobenzyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 1-cyclohexyl-4-(2-carboxyethyl)-7-(3-amidinobenzyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

1-ethyl-4-(2-carboxyethyl)-7-(4-amidinobutyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

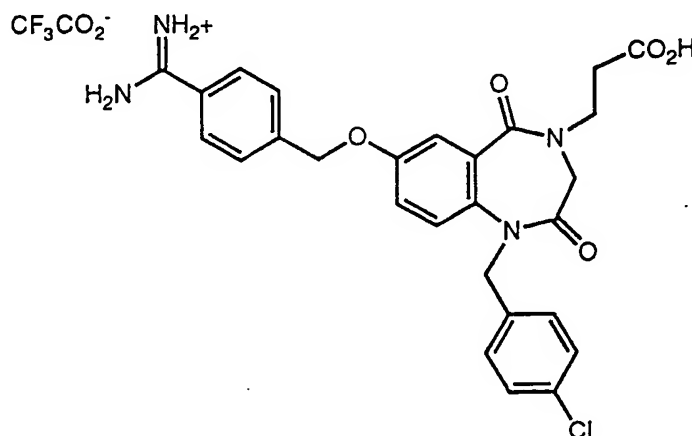
1-isopropyl-4-(2-carboxyethyl)-7-(4-amidinobutyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

5 1-isobutyl-4-(2-carboxyethyl)-7-(4-amidinobutyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,

1-cyclohexyl-4-(2-carboxyethyl)-7-(4-amidinobutyl)oxy-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate.

10

## Example 27



15 *1-(4-chlorobenzyl)-4-(2-carboxyethyl)-7-(4-amidinobenzyloxy)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate.*

15

a) To a stirred solution of 10.0 g (59.8 mmol) of 5-hydroxy-2-nitrobenzaldehyde and 12.3 g (62.8 mmol) of *p*-bromotoluenitrile in 50 ml of dry dimethyl-formamide was added 9.10 g (65.8 g) of anhydrous potassium carbonate. The mixture was warmed to 50 °C for 2 h and partitioned between water and ethyl acetate. The organic phase was washed with 1 N sodium bicarbonate, water, brine, and dried over anhydrous magnesium sulfate.

20

Concentration gave 13.8 g (82%) of 5-(4-cyanobenzyloxy)-2-nitrobenzaldehyde. <sup>1</sup>H NMR (300 MHz, d<sup>6</sup>-acetone) δ 10.39 (1H, s, Ar-CHO), 8.23 (1H, d, J=9.0 Hz, O-Ar-H), 7.85 (2H, d, J=8.3 Hz, NC-Ar-H<sub>2</sub>), 7.77 (2H, d, J=8.3 Hz, NC-Ar-H<sub>2</sub>), 7.48 (1H, dd, J=9.0, 2.9 Hz, O-Ar-H), 7.43 (1H, d, J=2.9 Hz, O-Ar-H), 5.50 (2H, s, ArCH<sub>2</sub>O).

25

b) Potassium permanganate (6.18 g, 39.1 mmol) was added in portions to a stirred solution of 13.8 g (48.9 mmol) of 5-(4-cyanobenzyloxy)-2-nitrobenzaldehyde and 1.58 g (4.89 mmol) of tetrabutylammonium bromide in 70 mL of pyridine. After 1 h, the solvent was removed in vacuo and the residue was partitioned between ethyl acetate and water. Sodium hydrogen

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sulfite was added to destroy excess permanganate and the aqueous phase was separated and extracted with ethyl acetate. The combined organics were washed with 1 N sodium hydrogen sulfate, brine, dried over anhydrous magnesium sulfate, and concentrated to give 14.6 g of 5-(4-cyanobenzyloxy)-2-nitrobenzoic acid as a powder, used without further purification. <sup>1</sup>H NMR (300 MHz, d<sup>6</sup>-acetone) δ 8.07 (1H, d, J=8.8 Hz, O-Ar-H), 7.86 (2H, d, J=8.3 Hz, NC-Ar-H<sub>2</sub>), 7.78 (2H, d, J=8.3 Hz, NC-Ar-H<sub>2</sub>), 7.39 (1H, d, J=2.9 Hz, O-Ar-H), 7.36 (1H, dd, J=8.8, 2.9 Hz, O-Ar-H), 5.48 (2H, s, ArCH<sub>2</sub>O).

c) Oxalyl chloride (5.70 mL, 65.4 mmol) was added slowly to a rapidly stirred suspension of 14.6 g (48.9 mmol) of 5-(4-cyanobenzyloxy)-2-nitrobenzoic acid and 2 mL of dimethylformamide in 100 mL of benzene. When gas evolution ceased, the solution was heated briefly to 60 °C, concentrated in vacuo, and redissolved in 75 mL of tetrahydrofuran. This solution was added dropwise to a rapidly stirred suspension of 8.50 g (55.3 mmol) of b-alanine ethyl ester hydrochloride and 21.1 g (251 mmol) of sodium bicarbonate in 325 mL of 2:1 tetrahydrofuran/water at 0 °C. The mixture was warmed to ambient temperature for 12 h and the tetrahydrofuran was removed in vacuo. The residue was partitioned between ethyl acetate and 1 N sodium bicarbonate and the aqueous phase was separated and extracted with ethyl acetate. The combined organics were washed with 1 N sodium hydrogen sulfate, water, brine, dried over anhydrous magnesium sulfate, and concentrated to give 17.7 g (91%) of N-(5-(4-cyanobenzyloxy)-2-nitrobenzoyl)-b-alanine ethyl ester as a tan powder. <sup>1</sup>H NMR (300 MHz, CDCl<sub>3</sub>) δ 8.11 (1H, d, J=8.3 Hz, O-Ar-H), 7.71 (2H, d, J=8.3 Hz, NC-Ar-H<sub>2</sub>), 7.52 (2H, d, J=8.3 Hz, NC-Ar-H<sub>2</sub>), 7.01 (2H, m, O-Ar-H<sub>2</sub>), 6.50 (1H, bt, J=5.9 Hz, N-H), 5.21 (2H, s, ArCH<sub>2</sub>O), 4.14 (2H, q, J=7.3 Hz, OCH<sub>2</sub>CH<sub>3</sub>), 3.69 (2H, bq, J=5.8 Hz, NCH<sub>2</sub>CH<sub>2</sub>), 2.70 (2H, t, J=5.8 Hz, NCH<sub>2</sub>CH<sub>2</sub>), 1.26 (3H, t, J=7.3 Hz, OCH<sub>2</sub>CH<sub>3</sub>); Exact mass (FAB, M+H<sup>+</sup>) calcd for C<sub>20</sub>H<sub>20</sub>N<sub>3</sub>O<sub>6</sub>: 398.1352; Found: 398.1343.

d) Tin(II) chloride (12.2 g, 64.4 mmol) was added to a solution of 5.12 g (12.9 mmol) of N-(5-(4-cyanobenzyloxy)-2-nitrobenzoyl)-b-alanine ethyl ester in 40 mL of 1:1 ethyl acetate/ethanol and the mixture heated to 70 °C for 15 min, cooled, and poured into a mixture of ice water, 1 N sodium bicarbonate, and ethyl acetate. 1 N sodium hydroxide was added carefully until a clear solution resulted and the aqueous phase was separated and extracted with ethyl acetate. The combined organics were washed with water, brine, dried over anhydrous magnesium sulfate, and concentrated to give a dark solid that was purified by chromatography (silica gel-60, 1:1 ethyl acetate/hexane) to give 2.51 g (53%) of N-(5-(4-cyanobenzyloxy)-2-aminobenzoyl)-b-alanine ethyl ester as a light yellow solid. <sup>1</sup>H NMR (300 MHz, CDCl<sub>3</sub>) δ 7.67 (2H, d, J=8.3 Hz, NC-Ar-H<sub>2</sub>), 7.52 (2H, d, J=8.3 Hz, NC-Ar-H<sub>2</sub>), 6.94 (1H, d, J=2.9 Hz, O-Ar-H), 6.89 (1H, dd, J=2.9, 8.8 Hz, O-Ar-H), 6.78 (1H, b, N-H), 6.66 (1H, d, J=8.8 Hz, O-Ar-H), 5.04 (2H, s, ArCH<sub>2</sub>O), 4.16 (2H, q, J=7.3 Hz, OCH<sub>2</sub>CH<sub>3</sub>), 3.66 (2H,



bq,  $J=5.8$  Hz,  $\text{NHCH}_2\text{CH}_2$ ), 2.62 (2H, t,  $J=5.8$  Hz,  $\text{NHCH}_2\text{CH}_2$ ), 1.27 (3H, t,  $J=7.3$  Hz,  $\text{OCH}_2\text{CH}_3$ ); Exact mass (FAB,  $\text{M}+\text{H}^+$ ) calcd for  $\text{C}_{20}\text{H}_{22}\text{N}_3\text{O}_4$ : 368.1610; Found: 368.1603.

e) A solution of 1.03 g (2.80 mmol) of N-(5-(4-cyanobenzoyloxy)-2-aminobenzoyl)-b-alanine ethyl ester, 394 mg (2.80 mmol) of p-chlorobenzaldehyde, and 20 mg of p-toluenesulfonic acid in 30 mL of toluene was heated at reflux through a Dean-Stark trap for 45 min. The solvent was removed in vacuo and the residue was dissolved in 15 mL of trifluoroacetic acid. Triethylsilane (1.34 mL, 8.40 mmol) was added and the mixture was stirred at room temperature for 12 h, concentrated in vacuo, and partitioned between 1 N aqueous sodium bicarbonate and ethyl acetate. The organic phase was washed with water, brine, dried over anhydrous magnesium sulfate, and concentrated. Chromatography (silica gel-60, 2:3 ethyl acetate/hexane) gave 820 mg (59%) of N-(5-(4-cyanobenzoyloxy)-2-(4-chlorobenzylamino)benzoyl)-b-alanine ethyl ester as a colorless solid.  $^1\text{H}$  NMR (300 MHz,  $\text{CDCl}_3$ )  $\delta$  7.66 (2H, d,  $J=7.8$  Hz, NC-Ar-H<sub>2</sub>), 7.51 (2H, d,  $J=7.8$  Hz, NC-Ar-H<sub>2</sub>), 7.27 (4H, s, ClArH<sub>4</sub>), 6.98 (1H, d,  $J=2.0$  Hz, O-Ar-H), 6.89 (1H, dd,  $J=2.0$ , 9.3 Hz, O-Ar-H), 6.78 (1H, b, N-H), 6.66 (1H, d,  $J=9.3$  Hz, O-Ar-H), 5.03 (2H, s, ArCH<sub>2</sub>O), 4.32 (2H, s, ClArCH<sub>2</sub>), 4.17 (2H, q,  $J=7.3$  Hz,  $\text{OCH}_2\text{CH}_3$ ), 3.65 (2H, bq,  $J=5.7$  Hz,  $\text{NHCH}_2\text{CH}_2$ ), 2.62 (2H, t,  $J=5.7$  Hz,  $\text{NHCH}_2\text{CH}_2$ ), 1.28 (3H, t,  $J=7.3$  Hz,  $\text{OCH}_2\text{CH}_3$ ); Exact mass (FAB,  $\text{M}+\text{H}^+$ ) calcd for  $\text{C}_{27}\text{H}_{27}\text{ClN}_3\text{O}_4$ : 492.1690; Found: 492.1681.

f) To a rapidly stirred mixture of 820 mg (1.67 mmol) of N-(5-(4-cyanobenzoyloxy)-2-(4-chlorobenzylamino)benzoyl)-b-alanine ethyl ester in 20 mL of 3:1 methylene chloride/water at ambient temperature was added 174 mL (2.00 mmol) of bromoacetyl bromide via syringe. After 6 h, an additional 100 mL (1.15 mmol) was added and the mixture stirred for 12 h. The aqueous phase was discarded and the organic layer washed with 1 N aqueous sodium bicarbonate, brine, dried over anhydrous magnesium sulfate, and concentrated. The residue was dissolved in 5 mL of dimethylformamide, 1.08 g (3.33 mmol) of cesium carbonate was added, and the mixture warmed briefly to 60 °C and allowed to stir at ambient temperature for 12 h. The resulting suspension was partitioned between ethyl acetate and water and the organic phase was separated and washed with water, brine, dried over anhydrous magnesium sulfate, and concentrated. Chromatography (silica gel-60, 1:1 ethyl acetate/hexane) gave 574 mg (65%) of 1-(4-chlorobenzyl)-4-(2-carboxyethyl)-7-(4-cyanobenzoyloxy)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester as a colorless foam.  $^1\text{H}$  NMR (300 MHz,  $\text{CDCl}_3$ )  $\delta$  7.68 (2H, d,  $J=8.3$  Hz, NC-Ar-H<sub>2</sub>), 7.51 (2H, d,  $J=8.3$  Hz, NC-Ar-H<sub>2</sub>), 7.35 (1H, d,  $J=3.0$  Hz, O-Ar-H), 7.24 (2H, d,  $J=9.7$  Hz, ClArH<sub>2</sub>), 7.10 (1H, d,  $J=8.8$  Hz, O-Ar-H), 7.07 (2H, d,  $J=9.7$  Hz, ClArH<sub>2</sub>), 7.01 (1H, dd,  $J=3.0$ , 8.8 Hz, O-Ar-H), 5.15 (1H, d,  $J=13.2$  Hz, ArCHHO), 5.10 (1H, d,  $J=13.2$  Hz, ArCHHO), 5.08 (1H, d,  $J=15.6$  Hz, ClArCH<sub>2</sub>), 5.91 (1H, d,  $J=15.6$  Hz, ClArCH<sub>2</sub>), 4.14 (2H, q,  $J=7.3$  Hz,  $\text{OCH}_2\text{CH}_3$ ), 4.13 (1H, d,  $J=14.7$  Hz, NCHHCON), 4.00 (1H, m, NCHHCH<sub>2</sub>), 3.89 (1H, d,  $J=14.7$  Hz, NCHHCON), 3.88

(NCHHCH<sub>2</sub>), 2.68 (2H, m, NHCH<sub>2</sub>CH<sub>2</sub>), 1.26 (3H, t, J=7.3 Hz, OCH<sub>2</sub>CH<sub>3</sub>); Exact mass (FAB, M+H<sup>+</sup>) calcd for C<sub>29</sub>H<sub>27</sub>ClN<sub>3</sub>O<sub>5</sub>: 532.1639; Found: 532.1628.

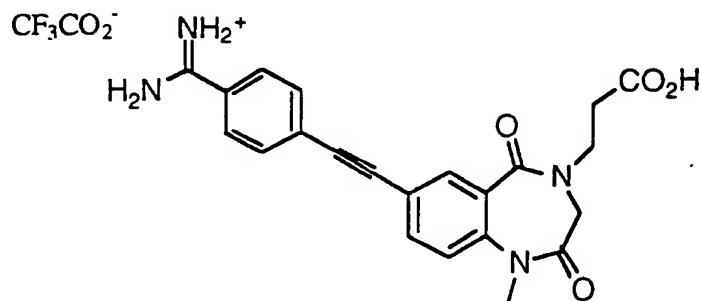
g) Hydrogen sulfide gas was bubbled into a solution of 295 mg (0.555 mmol) of 1-(4-chlorobenzyl)-4-(2-carboxyethyl)-7-(4-cyanobenzoyloxy)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester in 4 mL of 1:1 pyridine/triethylamine for 15 min and the mixture was warmed in a 50 °C oil bath for 12 h. After evaporation of the solvent, the residue was taken up in 4 mL of 3:1 methylene chloride/methyl iodide and heated at reflux for 1 h. The solution was again concentrated and the residue redissolved in 3 mL of methanol containing 0.5 g of ammonium acetate. The mixture was heated at 50 °C for 4 h, concentrated to a volume of 1 mL, and purified by reverse phase HPLC (Rainin Microsorb 1/2" C-18). The product was eluted with a solvent gradient of 0:100 acetonitrile (0.1% trifluoroacetic acid) /water (0.1% trifluoroacetic acid) to 50:50 acetonitrile (0.1% trifluoroacetic acid)/water (0.1% trifluoroacetic acid), time 0 to 30 min, flow=12 ml/min, R<sub>t</sub>=39.9 min, uv detection at 230 nM. Concentration of the clean fractions gave 215 mg (60%) of 1-(4-chlorobenzyl)-4-(2-carboxyethyl)-7-(4-amidinobenzoyloxy)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester trifluoroacetate as a colorless amorphous solid. <sup>1</sup>H NMR (300 MHz, d<sup>6</sup>-acetone) δ 8.60 (1H, b, NH), 7.95 (2H, d, J=8.3 Hz, HNC-Ar-H<sub>2</sub>), 7.73 (2H, d, J=8.3 Hz, HNC-Ar-H<sub>2</sub>), 7.40 (1H, d, J=9.3 Hz, O-Ar-H), 7.32-7.16 (6H, m, O-Ar-H, ArH<sub>4</sub>Cl), 5.34 (1H, d, J=16.1 Hz, NCHHAr), 5.27 (2H, s, ArCH<sub>2</sub>O), 4.94 (1H, d, J=16.1 Hz, NCHHAr), 4.24 (1H, d, J=14.7 Hz, NCHHCON), 4.11 (2H, q, J=7.3 Hz, OCH<sub>2</sub>CH<sub>3</sub>), 4.04 (1H, m, NCHHCH<sub>2</sub>), 3.94 (1H, d, J=14.7 Hz, NCHHCON), 3.80 (1H, m, NCHHCH<sub>2</sub>), 2.68 (2H, bt, J=6.8 Hz, NCH<sub>2</sub>CH<sub>2</sub>), 1.22 (3H, t, J=7.3 Hz, OCH<sub>2</sub>CH<sub>3</sub>); Exact mass (FAB, M+H<sup>+</sup>) calcd for C<sub>29</sub>H<sub>30</sub>ClN<sub>4</sub>O<sub>5</sub>: 549.1905; Found: 549.1912.

h) To a stirred, 0 °C solution of 104 mg (0.160 mmol) of 1-(4-chlorobenzyl)-4-(2-carboxyethyl)-7-(4-amidinobenzoyloxy)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester trifluoroacetate in 1.5 mL of tetrahydrofuran was added 0.48 mL of 1 N aqueous sodium hydroxide and the mixture warmed to ambient temperature over 2 h. The reaction was acidified with 0.10 mL of acetic acid, concentrated in vacuo, and purified by reverse phase HPLC (Rainin Microsorb 1/2" C-18). The product was eluted with a solvent gradient of 0:100 acetonitrile (0.1% trifluoroacetic acid) /water (0.1% trifluoroacetic acid) to 50:50 acetonitrile (0.1% trifluoroacetic acid)/water (0.1% trifluoroacetic acid), time 0 to 30 min, flow=12 ml/min, R<sub>t</sub>=29.1 min, uv detection at 230 nM. Lyophilization of the clean fractions gave 90.0 mg (90%) of 1-(4-chlorobenzyl)-4-(2-carboxyethyl)-7-(4-amidinobenzoyloxy)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate as a colorless powder. <sup>1</sup>H NMR (300 MHz, d<sup>6</sup>-acetone) δ 10.80 (1H, b, NH), 8.70 (1H, b, NH), 7.94 (2H, d, J=8.8 Hz, HNC-Ar-H<sub>2</sub>), 7.71 (2H, d, J=8.8 Hz, HNC-Ar-H<sub>2</sub>), 7.41 (1H, d, J=8.8 Hz, O-Ar-H), 7.32-7.16 (6H, m, O-Ar-H, ArH<sub>4</sub>Cl), 5.36 (1H, d, J=15.6 Hz, NCHHAr), 5.24 (2H, s, ArCH<sub>2</sub>O), 4.92 (1H, d, J=15.6

Hz, NCHHAr), 4.23 (1H, d, J=14.7 Hz, NCHHCON), 4.03 (1H, m, NCHHCH<sub>2</sub>), 3.98 (1H, d, J=14.7 Hz, NCHHCON), 3.81 (1H, m, NCHHCH<sub>2</sub>), 2.71 (2H, bt, J=6.8 Hz, NCH<sub>2</sub>CH<sub>2</sub>); Exact mass (FAB, M+H<sup>+</sup>) calcd for C<sub>27</sub>H<sub>26</sub>Cl N<sub>4</sub>O<sub>5</sub>: 521.1592; Found: 521.1581.

5

## Example 28



1-(methyl)-4-(2-carboxyethyl)-7-(4-amidinophenyl)ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate.

10

a) 1-methyl-4-(2-carboxyethyl)-7-(4-amidinophenyl)ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester was prepared using the method employed in part (h) of example 19. Thus, from 0.242 grams of 1-methyl-4-(2-carboxyethyl)-7-(4-cyanophenyl)ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester was prepared 0.145 grams (56%) of 1-methyl-4-(2-carboxyethyl)-7-(4-amidinophenyl)ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester. The product was purified by reverse phase HPLC (Column: Dynamax 60A C18, Method: 30:70 acetonitrile (0.1% trifluoroacetic acid) /water (0.1% trifluoroacetic acid), time 0 to 10 min, 30:70 acetonitrile (0.1% trifluoroacetic acid) /water (0.1% trifluoroacetic acid) to 70:30 acetonitrile (0.1% trifluoroacetic acid)/water (0.1% trifluoroacetic acid), time 10-40 min, 70:30 acetonitrile (0.1% trifluoroacetic acid)/water (0.1% trifluoroacetic acid), time 40-55min, flow=10 ml/min, R<sub>t</sub>=28.9 min, uv detection at 254 nm). <sup>1</sup>H NMR (300 MHz, D<sub>2</sub>O) δ 7.89 (1H, d, 3JHH=2Hz, C6-H), 7.80 (5H, m, C8-H, o,m-ArH), 7.46 (1H, d, 3JHH=9 Hz, C9-H), 4.17 (1H, d, J=15Hz, C3-H), 4.06 (1H, dt, 2JHH=14Hz, 3JHH=7Hz, NCHHCH<sub>2</sub>CO<sub>2</sub>), 3.92 (1H, d, J=15 Hz, C3-H), 3.77 (1H, dt, 2JHH=14Hz, 3JHH=7Hz, NCHHCH<sub>2</sub>CO<sub>2</sub>), 3.37 (3H, s, NCH<sub>3</sub>), 2.71 (2H, m, NCH<sub>2</sub>CH<sub>2</sub>CO<sub>2</sub>); Mass spectrum (FAB, M+H<sup>+</sup>) calcd for C<sub>24</sub>H<sub>25</sub>N<sub>4</sub>O<sub>4</sub>: 433.2; Found: 431.2.

Using the above procedure, but substituting the appropriate 4-(2-carboxyethyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester for 1-methyl-4-(2-carboxyethyl)-7-(4-cyanophenyl)ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester there was prepared, for example, the following compounds:

30

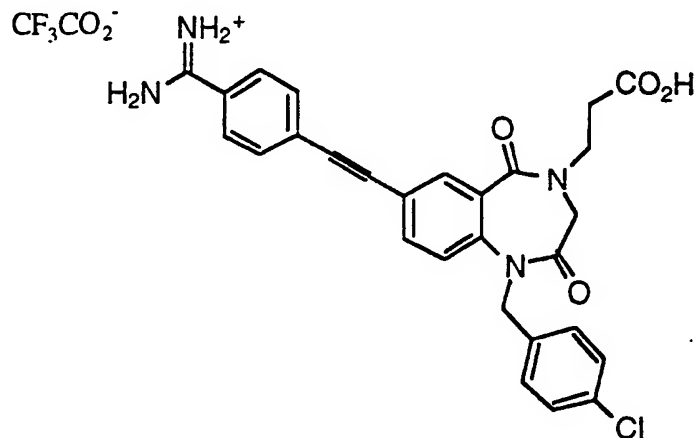
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- 1-methyl-4-(2-carboxyethyl)-7-(3-amidinophenyl)ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5,dione ethyl ester  
 1-ethyl-4-(2-carboxyethyl)-7-(4-amidinophenyl)ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5,dione ethyl ester  
 5 1-isopropyl-4-(2-carboxyethyl)-7-(4-amidinophenyl)ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5,dione ethyl ester  
 1-isobutyl-4-(2-carboxyethyl)-7-(4-amidinophenyl)ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5,dione ethyl ester  
 1-phenyl-4-(2-carboxyethyl)-7-(4-amidinophenyl)ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5,dione ethyl ester  
 10

- b) A magnetically stirred solution of 1-methyl-4-(2-carboxyethyl)-7-(4-amidinophenyl)ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl acetate (0.045 grams, 0.104 mmol) in 0.5mL methanol at room temperature was added 0.5mL 2N sodium hydroxide. After 30 mins, the reaction was quenched with 1mL acetic acid, concentrated *in vacuo*, diluted in a minimum of a 1:3 methanol:water mixture and purified by HPLC (Column: Dynamax 60A C18, Method: 10-50% MeCN/H<sub>2</sub>O 10ml/min. Detector: 254nm. Time(%MeCN): 0:00(10), 00:15(10, inj., detect on), 10:00(10), 40:00(50), 55:00(50), 59:58(10, detect off), 60:00(off), R<sub>f</sub>=32.3min) to yield 27mgs of 1-(methyl)-4-(2-carboxyethyl)-7-(4-amidinophenyl)ethynyl-3,4-dihydro-1H-1,4-benzodiazepinedione-2,5,dione trifluoroacetate. Exact mass (FAB, M+H<sup>+</sup>) calcd for C<sub>22</sub>H<sub>22</sub>N<sub>4</sub>O<sub>4</sub>: 405.1563, found: 405.1562  
 15  
 20

- Using the above procedure, but substituting the appropriate 4-(2-carboxyethyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester for 1-methyl-4-(2-carboxyethyl)-7-(4-amidinophenyl)ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5,dione ethyl ester there was prepared, for example, the following compounds:  
 1-methyl-4-(2-carboxyethyl)-7-(3-amidinophenyl)ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5,dione trifluoroacetate  
 1-ethyl-4-(2-carboxyethyl)-7-(4-amidinophenyl)ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5,dione trifluoroacetate  
 30 1-isopropyl-4-(2-carboxyethyl)-7-(4-amidinophenyl)ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5,dione trifluoroacetate  
 1-isobutyl-4-(2-carboxyethyl)-7-(4-amidinophenyl)ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5,dione trifluoroacetate  
 1-phenyl-4-(2-carboxyethyl)-7-(4-amidinophenyl)ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5,dione trifluoroacetate  
 35

## Example 29



5     1-(4-chlorobenzyl)-4-(2-carboxyethyl)-7-(4-amidinophenyl)ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate.

a) 1-(4-chlorobenzyl)-4-(2-carboxyethyl)-7-(4-cyanophenyl)ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester was prepared using the method described in part (f) of  
 10     example 1. Thus, from 0.265 grams (0.46mmol) of 1-(4-chlorobenzyl)-4-(2-carboxyethyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepinedione ethyl ester and 0.127 grams (4-cyanophenyl)ethyne was prepared 0.296 grams of 1-(4-chlorobenzyl)-4-(2-carboxyethyl)-7-(4-cyanophenyl)ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester (99%,  
 15     TLC, SiO<sub>2</sub>, 1:1 ethyl acetate/hexane, R<sub>f</sub>=0.38, uv positive). <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 7.99 (1H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, C6-H), 7.63 (2H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, ArH), 7.55 (2H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, ArH), 7.52 (1H, dd, <sup>4</sup>J<sub>HH</sub>=2Hz, <sup>3</sup>J<sub>HH</sub>=8Hz, C8-H), 7.24 (2H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, o-ArH NBn), 7.14 (1H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, C9-H), 7.02 (2H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, m-ArH NBn) 5.09 (1H, d, <sup>2</sup>J<sub>HH</sub>=15Hz, NCHHAr), 4.98 (1H, d, <sup>2</sup>J<sub>HH</sub>=15Hz, NCHHAr), 4.12 (4H, m, OCH<sub>2</sub>, C3-H, NCHHCH<sub>2</sub>CO<sub>2</sub>), 3.95 (1H, d, <sup>2</sup>J<sub>HH</sub>=15Hz, C3-H), 3.92 (1H, m, NCHHCH<sub>2</sub>CO<sub>2</sub>), 2.68 (2H,  
 20     m, CH<sub>2</sub>CO<sub>2</sub>), 1.24 (3H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, CH<sub>3</sub>)

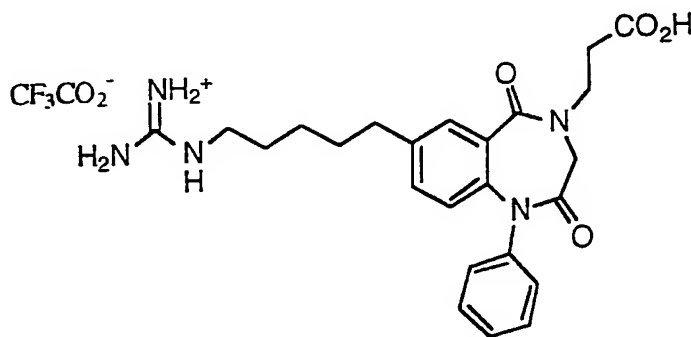
b) 1-(4-chlorobenzyl)-4-(2-carboxyethyl)-7-(4-amidinophenyl)ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester was prepared using the method described in part (h) of  
 25     example 19. Thus, from 0.15 grams (0.286mmol) of 1-(4-chlorobenzyl)-4-(2-carboxyethyl)-7-(4-cyanophenyl)ethynyl-3,4-dihydro-1H-1,4-benzodiazepinedione ethyl ester was prepared 0.03 grams (20%) of 1-(4-chlorobenzyl)-4-(2-carboxyethyl)-7-(4-amidinophenyl)ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester purified by HPLC (30-70%  
 MeOH/H<sub>2</sub>O 10ml/min. Detector: 254nm. Time(%MeOH): 0:00(30), 00:15(30, inj., detect on), 10:00(30), 40:00(70), 55:00(70), 59:58(30, detect off), 60:00(off), R<sub>f</sub>=38.0min). <sup>1</sup>H NMR  
 30     (CDCl<sub>3</sub>, dTMS) 7.99 (1H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, C6-H), 7.74 (2H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, ArH), 7.63 (2H, d,

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<sup>3</sup>J<sub>HH</sub>=8Hz, ArH), 7.54 (1H, dd, <sup>4</sup>J<sub>HH</sub>=2Hz, <sup>3</sup>J<sub>HH</sub>=8Hz, C8-H), 7.26 (2H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, *o*-ArH NBn), 7.16 (1H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, C9-H), 7.07 (2H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, *m*-ArH NBn) 5.10 (1H, d, <sup>2</sup>J<sub>HH</sub>=16Hz, NCHHAr), 4.98 (1H, d, <sup>2</sup>J<sub>HH</sub>=16Hz, NCHHAr), 4.13 (4H, m, OCH<sub>2</sub>, C3-H, NCHHCH<sub>2</sub>CO<sub>2</sub>), 3.95 (2H, m, C3-H, NCHHCH<sub>2</sub>CO<sub>2</sub>), 2.68 (2H, m, CH<sub>2</sub>CO<sub>2</sub>), 1.25 (3H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, CH<sub>3</sub>)

c) 1-(4-chlorobenzyl)-4-(2-carboxyethyl)-7-(4-amidinophenyl)etynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate was prepared using the method described in part (b) of example 27. Thus, 30mgs of 1-(4-chlorobenzyl)-4-(2-carboxyethyl)-7-(4-amidinophenyl)etynyl-3,4-dihydro-1H-1,4-benzodiazepinedione ethyl ester was saponified to yield 20mgs of 1-(4-chlorobenzyl)-4-(2-carboxyethyl)-7-(4-amidinophenyl)etynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate purified by HPLC (Column: Dynamax 60A C18, Method: 10-50% MeCN/H<sub>2</sub>O 10ml/min. Detector: 254nm. Time(%MeCN): 0:00(10), 00:15(10, inj., detect on), 10:00(10), 40:00(50), 55:00(50), 59:58(10, detect off), 60:00(off), R<sub>f</sub>=38.3min). <sup>1</sup>H NMR (D<sub>2</sub>O) 7.25 (6H, m, C6-H, ArH, C8-H), 6.98 (1H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, C9-H), 6.72 (2H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, *o*-ArH NBn), 6.57 (2H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, *m*-ArH NBn) 4.92 (1H, d, <sup>2</sup>J<sub>HH</sub>=16Hz, NCHHAr), NCHHAr overlaps with HOD, 3.70 (1H, d, <sup>2</sup>J<sub>HH</sub>=15Hz, C3-H), 3.54 (1H, m, NCHHCH<sub>2</sub>CO<sub>2</sub>), 3.43 (1H, d, <sup>2</sup>J<sub>HH</sub>=15Hz, C3-H), 3.30 (1H, m, NCHHCH<sub>2</sub>CO<sub>2</sub>), 2.59 (2H, m, CH<sub>2</sub>CO<sub>2</sub>). Exact mass (FAB, M+H<sup>+</sup>) calcd for C<sub>28</sub>H<sub>24</sub>N<sub>4</sub>O<sub>4</sub>Cl: 515.1514, found: 515.1486

## Example 30



25

1-phenyl-4-(2-carboxyethyl)-7-(5-guanidinopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate.

a) A solution of 2-chloro-5-nitrobenzoic acid (5.24 grams, 0.026mol), aniline (0.166mol, 15.5 grams), potassium carbonate (4.1 grams) and 0.1 grams copper oxide is heated refluxed for 2 hrs. The reaction mixture was allowed to cool to room temperature and diluted with water (100mL), and extracted with 1X50mL ether. The aqueous layer was made acidic by the

addition of conc. HCl and the mixture extracted 3X50mL ethyl acetate. The combined organics were dried over magnesium sulphate, filtered and concentrated *in vacuo*. The resulting residue, N-phenyl-5-nitroanthranilic acid (0.8 grams, 12%), was used without further purification. <sup>1</sup>H NMR (D<sub>2</sub>O, CD<sub>3</sub>OD) 8.84 (1H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, C6-H), 8.15 (1H, dd, <sup>2</sup>J<sub>HH</sub>=2Hz, <sup>3</sup>J<sub>HH</sub>=8Hz, C4-H), 7.42 (2H, d, *m*-NPh-H), 7.31 (3H, *m,o,p*-NPh-H), 7.15 (1H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, C3-H).

Using the above method, but substituting the appropriate substituted aniline for aniline there may be prepared, for example, the following compounds:

N-(4-methoxy)phenyl-5-nitroanthranilic acid, N-(4-chloro)phenyl-5-nitroanthranilic acid,  
10 N-(3-trifluoromethyl)phenyl-5-nitroanthranilic acid.

b) To a solution of N-phenyl-5-nitroanthranilic acid (0.8 grams, 3.1mmol), in 10mL methylene chloride and two drops dimethylformamide was added oxalyl chloride (0.3mL, 3.4mmol), and the reaction stirred for 30 mins. The mixture was concentrated in *vacuo*, diluted with  
15 15mL methylene chloride cooled to 0°C and allowed to react with b-alanine ethyl ester (3.4mmol, 0.522grams) and triethylamine (3.5mmol, 0.49mL). After 1 hour, the reaction was concentrated in *vacuo* and the resulting residue purified by column chromatography on SiO<sub>2</sub> using 25% ethyl acetate in hexane to 40% ethyl acetate in hexane (TLC, 1:2 ethyl acetate/hexane R<sub>f</sub>=0.31, uv positive) to yield 1 gram (90%) of N-(N-phenyl-5-nitroanthranoyl)-b-alanine ethyl ester. <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 10.38 (1H, bs, CO<sub>2</sub>H), 8.40 (1H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, C6-H anthranoyl), 8.08 (1H, dd, <sup>2</sup>J<sub>HH</sub>=2Hz, <sup>3</sup>J<sub>HH</sub>=9Hz, C4-H anthranoyl), 7.41 (2H, t, <sup>3</sup>J<sub>HH</sub>=8Hz, *m*-NPh-H), 7.24 (3H, m, *o,p*-NPh-H), 7.15 (1H, d, <sup>3</sup>J<sub>HH</sub>=9Hz, C3-H anthranoyl), 7.10 (1H, bt, <sup>3</sup>J<sub>HH</sub>=5Hz, C(O)NH), 4.21 (2H, q, <sup>3</sup>J<sub>HH</sub>=7Hz, OCH<sub>2</sub>), 3.75 (2H, q, <sup>3</sup>J<sub>HH</sub>=6Hz, NCH<sub>2</sub>CH<sub>2</sub>CO<sub>2</sub>), 2.71 (2H, t, <sup>3</sup>J<sub>HH</sub>=6Hz, CH<sub>2</sub>CO<sub>2</sub>), 1.31  
25 (3H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, CH<sub>3</sub>); <sup>13</sup>C NMR (CDCl<sub>3</sub>, dTMS) 172.5, 167.9, 151.9, 138.7, 137.1, 129.7, 128.1, 125.5, 124.6, 123.6, 114.8, 113.2, 61.1, 35.5, 33.7, 14.2.

Using the above method, but substituting the appropriate substituted N-aryl-5-nitroanthranilic acid for N-phenyl-5-nitroanthranilic acid there may be prepared, for example, the following compounds:

30 N-(N-(4-methoxy)phenyl-5-nitroanthranoyl)-b-alanine ethyl ester, N-(N-(4-chloro)phenyl-5-nitroanthranoyl)-b-alanine ethyl ester, N-(N-(4-trifluoromethyl)phenyl-5-nitroanthranoyl)-b-alanine ethyl ester.

c) A solution of 0.8 grams (2.3mmol) of N-(N-phenyl-5-nitroanthranoyl)-b-alanine ethyl ester in 5mL ethanol, 2.5mL 1N HCl, and 0.1 grams 10% palladium on carbon was placed under  
35 an atmosphere of hydrogen and stirred for 2 hrs. The mixture was filtered through Celite® washing with ethanol water mixture. The filtrate was concentrated *in vacuo* and the resulting residue diluted with 10mL water. The mixture was cooled to 0°C and allowed to react with sodium nitrite (158mgs, 2.3mmol) dissolved in 1mL water. After 2 hrs potassium

iodide (2.3 grams) in 3mL water was added and the reaction warmed to 50°C for 0.5 hrs. The cooled mixture was diluted with sat sodium bicarbonate and extracted 3X50mL ethyl acetate. The combined organics were dried over magnesium sulphate, filtered, concentrated *in vacuo*. and chromatographed on SiO<sub>2</sub>, using 20% ethyl acetate in hexane to 60% ethyl acetate and hexane (TLC, 25%ethyl acetate/hexane, R<sub>f</sub>=0.52 uv positive) to give 0.18 grams (18%) of N-(N<sup>2</sup>-phenyl-5-iodoanthranoyl)-b-alanine ethyl ester. <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 9.31 (1H, bs, NH), 7.66 (1H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, C6-H anthranoyl), 7.48 (1H, dd, <sup>2</sup>J<sub>HH</sub>=2Hz, <sup>3</sup>J<sub>HH</sub>=8Hz, C4-H anthranoyl), 7.32 (2H, m, NPh-H), 7.17 (2H, m, NPh-H), 7.10 (1H, d, <sup>3</sup>J<sub>HH</sub>=9Hz, C3-H anthranoyl), 7.04 (1H, m, NPh-H), 6.85 (1H, bt, <sup>3</sup>J<sub>HH</sub>=5Hz, C(O)NH), 4.20 (2H, q, <sup>3</sup>J<sub>HH</sub>=7Hz, OCH<sub>2</sub>), 3.69 (2H, q, <sup>3</sup>J<sub>HH</sub>=6Hz, NCH<sub>2</sub>CH<sub>2</sub>CO<sub>2</sub>), 2.65 (2H, t, <sup>3</sup>J<sub>HH</sub>=6Hz, CH<sub>2</sub>CO<sub>2</sub>), 1.29 (3H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, CH<sub>3</sub>); <sup>13</sup>C NMR (CDCl<sub>3</sub>, dTMS) 172.7, 168.0, 151.2, 140.7, 140.5, 135.9, 129.4, 123.0, 121.2, 120.2, 117.2, 77.9, 60.9, 35.3, 33.9, 14.2.

Using the above method, but substituting the appropriate substituted N-(N-aryl-5-nitroanthranoyl)-b-alanine ethyl ester for N-(N-phenyl-5-nitroanthranoyl)-b-alanine ethyl ester there may be prepared, for example, the following compounds:

15 N-(N-(4-methoxy)phenyl-5-iodoanthranoyl)-b-alanine ethyl ester, N-(N-(4-chloro)phenyl-5-iodoanthranoyl)-b-alanine ethyl ester, N-(N-(4-trifluoromethyl)phenyl-5-iodoanthranoyl)-b-alanine ethyl ester.

20 d) To a magnetically stirred biphasic mixture of N-(N-phenyl-5-iodoanthranoyl)-b-alanine ethyl ester (180mgs, 0.41mmol), 10mL methylene chloride, 10mL water was added bromoacetyl bromide (0.054mL, 0.45mmol). After 2 hrs, the layers were separated. The organic layer was dried over sodium sulphate, decanted, and concentrated in vacuo. The resulting residue was dissolved in dimethylformamide (3mL) and cesium carbonate (2.5 molar excess, 1.0mmol) added. After 30 mins, the mixture was diluted with 75mL water and extracted with 3X50mL ethyl acetate. The combined organics were dried over sodium sulphate, decanted, and concentrated in vacuo. The resulting residue was further purified by column chromatography on SiO<sub>2</sub>, using 35% ethyl acetate in hexane to 65% ethyl acetate and hexane to yield 0.142 grams (72%) of 1-phenyl-4-(2-carboxyethyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester (TLC, 1:1 ethyl acetate/hexane, R<sub>f</sub>=0.29, uv positive). <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 8.20 (1H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, C6-H), 7.60 (1H, dd, <sup>2</sup>J<sub>HH</sub>=2Hz, <sup>3</sup>J<sub>HH</sub>=8Hz, C8-H), 7.38 (3H, m, NPh-H), 7.20 (2H, m, NPh-H), 6.53 (1H, d, <sup>3</sup>J<sub>HH</sub>=9Hz, C9-H), 4.21 (1H, d, <sup>2</sup>J<sub>HH</sub>=15Hz, C3-H), 4.11 (2H, q, <sup>3</sup>J<sub>HH</sub>=7Hz, OCH<sub>2</sub>), 4.01 (1H, d, <sup>2</sup>J<sub>HH</sub>=15Hz, C3-H), 3.95 (2H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, NCH<sub>2</sub>CH<sub>2</sub>CO<sub>2</sub>), 2.72 (2H, m, CH<sub>2</sub>CO<sub>2</sub>), 1.24 (3H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, CH<sub>3</sub>); <sup>13</sup>C NMR (CDCl<sub>3</sub>, dTMS) 171.2, 167.2, 165.6, 140.7, 140.5, 139.8, 139.3, 130.7, 129.5, 128.1, 128.0, 126.1, 89.9, 60.8, 52.6, 45.2, 32.6, 14.1.

Using the above method, but substituting the appropriate substituted N-aryl-5-iodoanthranilic acid for N-(N-phenyl-5-iodoanthranoyl)-b-alanine ethyl ester there may be prepared, for example, the following compounds:



1-(4-methoxy)phenyl-4-(2-carboxyethyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester, 1-(4-chloro)phenyl-4-(2-carboxyethyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester, 1-(4-trifluoromethyl)phenyl-4-(2-carboxyethyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester.

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e) 1-phenyl-4-(2-carboxyethyl)-7-(N-Boc-5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester was prepared using the method described in part (f) of example 1. Thus, 71mgs of 1-phenyl-4-(2-carboxyethyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepinedione-2,5,dione ethyl ester and 30mgs of N-Boc-5-amino-1-pentyne yielded 10 61mgs (76%) of 1-phenyl-4-(2-carboxyethyl)-7-(N-Boc-5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester. <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 7.90 (1H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, C6-H), 7.37 (6H, m, C8-H, NPh-H), 6.69 (1H, d, <sup>3</sup>J<sub>HH</sub>=9Hz, C9-H), 4.72 (1H, bs, NH), 4.21 (1H, d, <sup>2</sup>J<sub>HH</sub>=15Hz, C3-H), 4.11 (2H, q, <sup>3</sup>J<sub>HH</sub>=7Hz, OCH<sub>2</sub>), 4.00 (1H, d, <sup>2</sup>J<sub>HH</sub>=15Hz, C3-H), 3.96 (2H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, NCH<sub>2</sub>CH<sub>2</sub>CO<sub>2</sub>), 3.25 (2H, bq, <sup>3</sup>J<sub>HH</sub>=7Hz, CH<sub>2</sub>NH), 2.72 (2H, m, CH<sub>2</sub>CO<sub>2</sub>), 2.44 (2H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, CCCH<sub>2</sub>), 1.76 (2H, p, <sup>3</sup>J<sub>HH</sub>=7Hz, CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>), 1.43 (9H, s, C(CH<sub>3</sub>)<sub>3</sub>), 1.23 (3H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, CH<sub>3</sub>); <sup>13</sup>C NMR (CDCl<sub>3</sub>, dTMS) 171.3, 167.3, 166.4, 155.9, 140.0, 139.7, 134.5, 133.8, 129.4, 129.1, 128.1, 127.9, 124.3, 121.7, 90.0, 79.5, 60.8, 52.6, 45.1, 39.7, 32.7, 28.7, 28.4, 16.9, 14.1.

Using the above procedure, but substituting the appropriate 1-aryl-4-(2-carboxyethyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepinedione-2,5,dione ethyl ester for 1-phenyl-4-(2-carboxyethyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepinedione-2,5,dione ethyl ester and amino or cyano alkyne for N-Boc-5-amino-1-pentyne there may be prepared, for example, the following compounds:

1-(4-methoxy)phenyl-4-(2-carboxyethyl)-7-(N-Boc-5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,

1-(4-chloro)phenyl-4-(2-carboxyethyl)-7-(N-Boc-5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester ,

1-(3-trifluorophenyl)phenyl-4-(2-carboxyethyl)-7-(N-Boc-5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester

30 1-(4-methoxy)phenyl-4-(2-carboxyethyl)-7-(N-Boc-6-amino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,

1-(4-chloro)phenyl-4-(2-carboxyethyl)-7-(N-Boc-6-amino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester ,

1-(3-trifluorophenyl)phenyl-4-(2-carboxyethyl)-7-(N-Boc-6-amino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester

35 1-(4-methoxy)phenyl-4-(2-carboxyethyl)-7-(4-cyanophenyl)ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,

1-(4-chloro)phenyl-4-(2-carboxyethyl)-7-(4-cyanophenyl)ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester ,

1-(3-trifluorophenyl)phenyl-4-(2-carboxyethyl)-7-(4-cyanophenyl)ethynyl-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester

- f) Using the method described in part (a) of example 3 1-phenyl-4-(2-carboxyethyl)-7-(N-Boc-5-aminopentyl)-3,4-dihydro-1H-1,4-benzodiazepinedione-2,5-dione ethyl ester was prepared from 1-phenyl-4-(2-carboxyethyl)-7-(N-Boc-5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester in quantitative yield. <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 7.67 (1H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, C6-H), 7.40 (5H, m, NPh-H), 7.13 (1H, dd, <sup>4</sup>J<sub>HH</sub>=2Hz, <sup>3</sup>J<sub>HH</sub>=8Hz, C8-H), 6.70 (1H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, C9-H), 4.55 (1H, bs, NH), 4.24 (1H, d, <sup>2</sup>J<sub>HH</sub>=15Hz, C3-H), 4.11 (2H, q, <sup>3</sup>J<sub>HH</sub>=7Hz, OCH<sub>2</sub>), 3.97 (2H, m, C3-H, NCH<sub>2</sub>CH<sub>2</sub>CO<sub>2</sub>), 3.10 (2H, bq, <sup>3</sup>J<sub>HH</sub>=7Hz, CH<sub>2</sub>NH), 2.72 (2H, m, CH<sub>2</sub>CO<sub>2</sub>), 2.60 (2H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, ArCH<sub>2</sub>), 1.62 (2H, p, <sup>3</sup>J<sub>HH</sub>=7Hz), 1.47 (2H, p, <sup>3</sup>J<sub>HH</sub>=7Hz), 1.41 (9H, s, C(CH<sub>3</sub>)<sub>3</sub>), 1.34 (2H, m), 1.23 (3H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, CH<sub>3</sub>); <sup>13</sup>C NMR (CDCl<sub>3</sub>, dTMS) 171.3, 167.6, 167.2, 155.9, 140.4, 140.3, 138.6, 132.0, 130.1, 129.3, 128.9, 128.1, 127.6, 124.3, 60.7, 52.7, 45.1, 40.4, 34.9, 32.7, 30.6, 29.9, 28.4, 26.4, 14.1.

- Using the above procedure, but substituting the appropriate 1-aryl-4-(2-carboxyethyl)-7-alkynyl-3,4-dihydro-1H-1,4-benzodiazepinedione-2,5-dione ethyl ester for 1-phenyl-4-(2-carboxyethyl)-7-(N-Boc-5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester there may be prepared, for example, the following compounds:

- 1-(4-methoxy)phenyl-4-(2-carboxyethyl)-7-(N-Boc-5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,  
1-(4-chloro)phenyl-4-(2-carboxyethyl)-7-(N-Boc-5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester ,  
1-(3-trifluorophenyl)phenyl-4-(2-carboxyethyl)-7-(N-Boc-5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester  
1-(4-methoxy)phenyl-4-(2-carboxyethyl)-7-(N-Boc-6-amino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,  
1-(4-chloro)phenyl-4-(2-carboxyethyl)-7-(N-Boc-6-amino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester ,  
1-(3-trifluorophenyl)phenyl-4-(2-carboxyethyl)-7-(N-Boc-6-amino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester

- g) To a solution of 13.3mgs (0.025mmol) 1-phenyl-4-(2-carboxyethyl)-7-(N-Boc-5-aminopentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester in 1mL ethyl acetate and 0.5mL triethylsilane was added 2mL sat. solution of HCl in ethyl acetate. After 30 mins, the reaction mixture was concentrated *in vacuo*, diluted in a minimum of THF and 18mL of a 30% by weight hydrogen peroxide. An equivalent of lithium hydroxide was added and the reaction stirred for 1 hrs. The reaction was quenched with sodium sulfite and

concentrated to yield a residue, which was diluted in a methanol water mixture and purified by HPLC (10-50% MeCN/H<sub>2</sub>O 10ml/min. Detector: 254nm. Time(%MeCN): 0:00(10), 00:15(10, inj., detect on), 10:00(10), 40:00(50), 55:00(50), 59:58(10, detect off), 60:00(off), R<sub>f</sub>=33.5min) to yield 1-phenyl-4-(2-carboxyethyl)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-

- 5 1,4-benzodiazepine-2,5-dione (6mgs). <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 7.12 (1H, bs, C<sub>6</sub>-H), 6.97 (3H, m, *m,p*-NPh-H), 6.79 (1H, bd, <sup>3</sup>J<sub>HH</sub>=8Hz, C<sub>8</sub>-H), 6.74 (2H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, *o*-NPh-H), 6.36 (1H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, C<sub>9</sub>-H), 3.89 (1H, d, <sup>2</sup>J<sub>HH</sub>=15Hz, C<sub>3</sub>-H), 3.74 (1H, dt, <sup>2</sup>J<sub>HH</sub>=14Hz, <sup>3</sup>J<sub>HH</sub>=7Hz, NCHHCH<sub>2</sub>CO<sub>2</sub>), 3.47 (1H, d, <sup>2</sup>J<sub>HH</sub>=15Hz, C<sub>3</sub>-H), 3.21 (1H, dt, <sup>2</sup>J<sub>HH</sub>=14Hz, <sup>3</sup>J<sub>HH</sub>=7Hz, NCHHCH<sub>2</sub>CO<sub>2</sub>), 2.43 (2H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, CH<sub>2</sub>NH<sub>2</sub>), 2.25 (2H, m, CH<sub>2</sub>CO<sub>2</sub>),  
10 2.15 (2H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, ArCH<sub>2</sub>), 1.12 (4H, m, ), 0.83 (2H, m)

Using the above procedure, but substituting the appropriate 1-aryl-4-(2-carboxyethyl)-7-alkynyl-3,4-dihydro-1H-1,4-benzodiazepinedione-2,5,dione ethyl ester for 1-phenyl-4-(2-carboxyethyl)-7-(N-Boc-5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester there may be prepared, for example, the following

- 15 compounds:  
1-(4-methoxy)phenyl-4-(2-carboxyethyl)-7-(N-Boc-5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,  
1-(4-chloro)phenyl-4-(2-carboxyethyl)-7-(N-Boc-5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester ,  
20 1-(3-trifluorophenyl)phenyl-4-(2-carboxyethyl)-7-(N-Boc-5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester  
1-(4-methoxy)phenyl-4-(2-carboxyethyl)-7-(N-Boc-6-amino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester,  
1-(4-chloro)phenyl-4-(2-carboxyethyl)-7-(N-Boc-6-amino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester ,  
25 1-(3-trifluorophenyl)phenyl-4-(2-carboxyethyl)-7-(N-Boc-6-amino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester

- h) 6mgs of 1-phenyl-4-(2-carboxyethyl)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5,dione was dissolved in sat 5% potassium carbonate and an excess of formamidine sulphonic acid (15mgs). After 30mins, the reaction was quenched with acetic acid , concentrated *in vacuo*, diluted in a minimum of a water/methanol mixture and purified by HPLC (Column: Dynamax 60A C18, Method: 10-50% MeCN/H<sub>2</sub>O 10ml/min. Detector: 254nm. Time(%MeCN): 0:00(10), 00:15(10, inj., detect on), 10:00(10), 40:00(50), 55:00(50),  
35 59:58(10, detect off), 60:00(off), R<sub>f</sub>=37.1min). <sup>1</sup>H NMR (D<sub>2</sub>O) 7.08 (1H, bs, ArH C<sub>6</sub>-H), 6.93 (3H, m), 6.63 (1H, bd, <sup>3</sup>J<sub>HH</sub>=4Hz), 6.21 (2H, d, <sup>3</sup>J<sub>HH</sub>=8Hz), 3.79 (1H, d, <sup>2</sup>J<sub>HH</sub>=15Hz, C<sub>3</sub>-H), 3.74 (1H, dt, <sup>2</sup>J<sub>HH</sub>=14Hz, <sup>3</sup>J<sub>HH</sub>=7Hz, NCHHCH<sub>2</sub>CO<sub>2</sub>), 3.42 (1H, d, <sup>2</sup>J<sub>HH</sub>=15Hz, C<sub>3</sub>-H), 3.19 (1H, dt, <sup>2</sup>J<sub>HH</sub>=14Hz, <sup>3</sup>J<sub>HH</sub>=7Hz, NCHHCH<sub>2</sub>CO<sub>2</sub>), 2.41 (2H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, CH<sub>2</sub>NH),

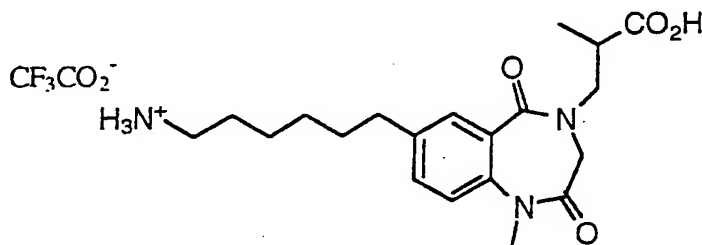
2.23 (2H, m, CH<sub>2</sub>CO<sub>2</sub>), 2.12 (2H, bt, <sup>3</sup>J<sub>HH</sub>=7Hz), 1.02 (4Hz, m), 0.79 (2H, m) Exact mass (FAB, M+H<sup>+</sup>) calculated for C<sub>24</sub>H<sub>30</sub>N<sub>5</sub>O<sub>4</sub>; 452.2298, found 452.2270.

Using the above procedure, but substituting the appropriate 1-aryl-4-(2-carboxyethyl)-7-alkyl-3,4-dihydro-1H-1,4-benzodiazepinedione-2,5-dione ethyl ester for 1-phenyl-4-(2-carboxyethyl)-7-(N-Boc-5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione ethyl ester there may be prepared, for example, the following compounds:

- 1-(4-methoxy)phenyl-4-(2-carboxyethyl)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 10 1-(4-chloro)phenyl-4-(2-carboxyethyl)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate ,
- 1-(3-trifluorophenyl)phenyl-4-(2-carboxyethyl)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate
- 1-(4-methoxy)phenyl-4-(2-carboxyethyl)-7-(6-amino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate,
- 15 1-(4-chloro)phenyl-4-(2-carboxyethyl)-7-(6-amino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate ,
- 1-(3-trifluorophenyl)phenyl-4-(2-carboxyethyl)-7-(6-amino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate

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## Example 31



- 25 (±)-1-methyl-4-(2-carboxy-2-methylethyl)-7-(6-amino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate.

a) In a 1L beaker was placed 40 grams (0.267mmol) of N-methylanthranilic acid dissolved in 300mL water and 26.7mL conc. HCl cooled to 20°C. A solution of ICL, prepared by dissolving 44grams of ICl to a cooled (0°C) solution of 167mL water and 45ml conc. HCl, was rapidly added to the stirred N-methylanthranilic acid solution. The reaction was allowed to stir for 2hrs, filtered on a medium glass frit funnel, washing the solids with water. The green powder isolated was dried under vacuum to yield 70 grams (96%) of 5-iodo-N-methylanthranilic acid. <sup>1</sup>H NMR (DMSO-d<sub>6</sub>) 7.61 (1H, bs, C6-H), 7.21 (1H, bd, <sup>3</sup>J<sub>HH</sub>=8Hz,

C4-H), 6.19 (1H, d,  $^3J_{HH}=8\text{Hz}$ , C3-H), 2.43 (3H, s, NCH<sub>3</sub>);  $^{13}\text{C}$  NMR (DMSO- $d_6$ ) 169.0, 151.4, 142.7, 139.6, 114.2, 112.7, 74.2, 29.6.

b) To a mechanically stirred solution of 5-iodo-N-methylantranilic acid (40 grams 0.144 mol), potassium carbonate (19.95 grams, 0.144mol), and 500ml water at 0°C was added phosgene (1.93M solution in toluene, 0.2mol, 104mL) over a period of 1hr. the resulting slurry was stirred for an additional 3 hrs and the solids isolated by filtration, washing with water, and a small amount (70mL) of ether. The solids were dried overnight *in vacuo* to yield 34 grams (78%) of 5-iodo-N-methylisatoic anhydride.  $^1\text{H}$  NMR (DMSO- $d_6$ ) 8.18 (1H, bs, C6-H), 8.08 (1H, bd,  $^3J_{HH}=8\text{Hz}$ , C4-H), 7.12 (1H, d,  $^3J_{HH}=8\text{Hz}$ , C3-H), 3.42 (3H, s, NCH<sub>3</sub>).

c) A solution of ( $\pm$ )-3-amino-2-methylpropionic acid (1.9 grams, 18.77mmol), *p*-toluenesulphonic acid, in 50mL methanol was heated to reflux for 62hrs. the reaction was allowed to cool to room temperature, concentrated *in vacuo* to yield an oil. 1/2 by volume of this oil was dissolved in 10 mL dimethylformamide and allowed to react with 5-iodo-N-methylisatoic anhydride (2.5 grams, 8.4mmol) and 7.0mL triethylamine (5-fold excess) and a catalytic amount of dimethylaminopyridine (50mgs). the reaction was heated to 50°C for 1hr. The reaction was allowed to cool to room temperature, diluted with 100mL water and extracted 3X50mL ethyl acetate. The combined organics were dried over sodium sulphate, decanted, and concentrated *in vacuo*. The resulting residue was further purified by column chromatography (SiO<sub>2</sub>, 25%ethyl acetate in hexane to 50% ethyl acetate and hexane) to yield 1.5 grams (47%) of ( $\pm$ )-N-(N<sup>2</sup>-methyl-2-amino-5-iodobenzoyl)-3-amino-2-methylpropionate methyl ester.  $^1\text{H}$  NMR (CDCl<sub>3</sub>, dTMS) 7.50 (1H, d,  $^4J_{HH}=2\text{Hz}$ , C6-H), 7.43 (1H, dd,  $^4J_{HH}=2\text{Hz}$ ,  $^3J_{HH}=9\text{Hz}$ , C4-H), 6.79 (1H, bt, NHCH<sub>2</sub>) 6.35 (1H, d,  $^3J_{HH}=9\text{Hz}$ , C3-H), 3.64 (3H, s, OCH<sub>3</sub>), 3.51 (1H, ddd,  $^2J_{HH}=14\text{Hz}$ ,  $^3J_{HH}=6\text{Hz}$ ,  $^3J_{HH}=5\text{Hz}$ , NHCH<sub>2</sub>CH), 3.40 (1H, ddd,  $^2J_{HH}=14\text{Hz}$ ,  $^3J_{HH}=8\text{Hz}$ ,  $^3J_{HH}=6\text{Hz}$ , NHCH<sub>2</sub>CH), 2.75 (4H, m, CHCO<sub>2</sub>, NCH<sub>3</sub>), 1.15 (3H, d,  $^3J_{HH}=7\text{Hz}$ , CHCH<sub>3</sub>);  $^{13}\text{C}$  NMR 175.9, 168.7, 149.9, 141.0, 135.7, 117.6, 113.4, 74.2, 52.1, 42.1, 39.6, 29.7, 15.1

d) To a biphasic mixture of ( $\pm$ )-N-(N-methyl-2-amino-5-iodobenzoyl)-3-amino-2-methylpropionate methyl ester (2.0mmol, 0.752grams), 8mL methylene chloride, and 8mL water was added bromoacetyl bromide (0.174mL, 1.5 molar equivalents) and the reaction stirred for 2hrs. The layers were separated and the organics were dried over sodium sulphate, decanted, and concentrated *in vacuo*. The resulting residue was diluted in 10mL dimethylformamide and a 3-fold excess of cesium carbonate (6.0mmol, 1.95 grams). After 30 mins the mixture was diluted with 75mL water and extracted 2X50mL ethyl acetate. The combined organics were dried over magnesium sulphate, filtered, and concentrated *in vacuo*. The resulting residue was further purified by column chromatography (SiO<sub>2</sub>, 50% ethyl acetate/hexane to 70% ethyl acetate/hexane, TLC,  $R_f=0.21$ , uv positive) to yield ( $\pm$ )-1-

methyl-4-(2-carboxy-2-methylethyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester (0.5 grams, 61%, mp=138-140°C). <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 8.13 (1H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, C6-H), 7.79 (1H, dd, <sup>4</sup>J<sub>HH</sub>=2Hz, <sup>3</sup>J<sub>HH</sub>=9Hz, C8-H), 6.92 (1H, d, <sup>3</sup>J<sub>HH</sub>=9Hz, C9-H), 3.80 (7H, s, NCH<sub>2</sub>CHCO<sub>2</sub>, C3-H, OCH<sub>3</sub>), 3.34 (3H, two s, NCH<sub>3</sub>, diastereotopic resonances are observed), 1.19 (1.5H, d, <sup>3</sup>J<sub>HH</sub>=7Hz, CHCH<sub>3</sub>), 1.16 (1.5H, d, <sup>3</sup>J<sub>HH</sub>=7Hz, CHCH<sub>3</sub>).

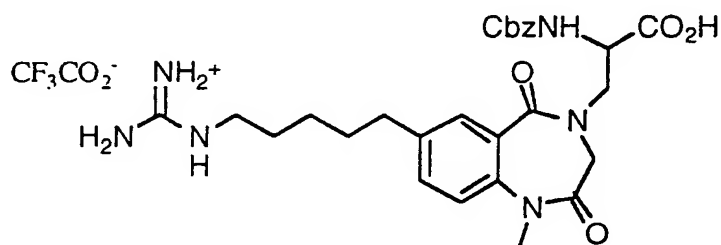
e) (±)-1-methyl-4-(2-carboxy-2-methylethyl)-7-(N-Boc-5-amino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester was prepared using the method described in part (f) of example 1. Thus, 0.15 grams (0.36mmol) of (±)-1-methyl-4-(2-carboxy-2-methylethyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepinedione-2,5-dione methyl ester and 0.11 grams (0.54mmol) of N-Boc-6-amino-1-hexyne yielded 0.164 grams (94%) of (±)-1-methyl-4-(2-carboxy-2-methylethyl)-7-(N-Boc-5-amino-1-hexynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester (TLC, 1:1 ethyl acetate/hexane, R<sub>f</sub>=0.23). <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 7.79 (1H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, C6-H), 7.43 (1H, dd, <sup>4</sup>J<sub>HH</sub>=2Hz, <sup>3</sup>J<sub>HH</sub>=8Hz, C8-H), 7.06 (1H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, C9-H), 4.67 (1H, bs, NHBoc), 4.04 (0.5H, d, <sup>2</sup>J<sub>HH</sub>=15Hz, C3-H), 3.94 (0.5H, d, <sup>2</sup>J<sub>HH</sub>=15Hz, C3-H), 3.93 (0.5H, d, <sup>2</sup>J<sub>HH</sub>=15Hz, C3-H), 3.73 (2.5H, m, NCH<sub>2</sub>CH, C3-H), 3.66 (1.5H, s, OCH<sub>3</sub>), 3.59 (1.5H, s, OCH<sub>3</sub>), 3.30 (3H, bs, NCH<sub>3</sub>), 3.10 (2H, bq, CH<sub>2</sub>NH), 2.93 (1H, m, CHCO<sub>2</sub>), 2.38 (2H, bt, <sup>3</sup>J<sub>HH</sub>=7Hz, CCCH<sub>2</sub>), 1.58 (4H, m), 1.38 (9H, s, C(CH<sub>3</sub>)<sub>3</sub>), 1.14 (1.5H, d, <sup>3</sup>J<sub>HH</sub>=7Hz, CHCH<sub>3</sub>), 1.11 (1.5H, d, <sup>3</sup>J<sub>HH</sub>=7Hz, CHCH<sub>3</sub>); <sup>13</sup>C NMR 175.1, 174.6, 168.7, 166.9, 156.1, 139.9, 134.9, 134.0, 128.7, 121.8, 120.9, 91.5, 79.3, 79.1, 52.5, 52.1, 52.0, 51.5, 40.2, 38.5, 38.4, 34.9, 34.8, 29.4, 28.5, 25.8, 19.2, 16.1, 14.6.

f) (±)-1-methyl-4-(2-carboxy-2-methylethyl)-7-(N-Boc-5-amino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester was prepared using the method described in part (a) of example 3. Thus, 0.085 grams (0.17mmol) of (±)-1-methyl-4-(2-carboxy-2-methylethyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepinedione-2,5-dione methyl ester yielded 0.081 grams (95%) of (±)-1-methyl-4-(2-carboxy-2-methylethyl)-7-(N-Boc-5-amino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester (TLC, 3:1 ethyl acetate/hexane, R<sub>f</sub>=0.62). <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 7.58 (1H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, C6-H), 7.27 (1H, dd, <sup>4</sup>J<sub>HH</sub>=2Hz, <sup>3</sup>J<sub>HH</sub>=8Hz, C8-H), 7.06 (1H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, C9-H), 4.60 (1H, bs, NHBoc), 4.07 (0.5H, d, <sup>2</sup>J<sub>HH</sub>=15Hz, C3-H), 3.98 (0.5H, d, <sup>2</sup>J<sub>HH</sub>=15Hz, C3-H), 3.97 (0.5H, d, <sup>2</sup>J<sub>HH</sub>=15Hz, C3-H), 3.75 (2.5H, m, NCH<sub>2</sub>CH, C3-H), 3.67 (1.5H, s, OCH<sub>3</sub>), 3.60 (1.5H, s, OCH<sub>3</sub>), 3.33 (3H, two s, NCH<sub>3</sub>), 3.06 (2H, bq, CH<sub>2</sub>NH), 2.95 (1H, m, CHCO<sub>2</sub>), 2.58 (2H, bt, <sup>3</sup>J<sub>HH</sub>=7Hz, ArCH<sub>2</sub>), 1.59 (2H, m), 1.40 (11H, m, C(CH<sub>3</sub>)<sub>3</sub>), 1.30 (4H, m), 1.15 (1.5H, d, <sup>3</sup>J<sub>HH</sub>=7Hz, CHCH<sub>3</sub>), 1.13 (1.5H, d, <sup>3</sup>J<sub>HH</sub>=7Hz, CHCH<sub>3</sub>); <sup>13</sup>C NMR 174.9, 174.5, 168.8, 167.4, 155.9, 140.2, 138.6, 132.1, 130.1, 128.2, 120.7, 78.8, 52.4, 52.0, 51.9, 51.2, 40.3, 38.4, 34.8, 34.7, 30.8, 29.8, 28.7, 28.3, 26.4, 14.9, 14.4.

( $\pm$ )-1-methyl-4-(2-carboxy-2-methylethyl)-7-(6-amino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate was prepared using the method described in part (g) of example 1. Thus 0.04grams of ( $\pm$ )-1-methyl-4-(2-carboxy-2-methylethyl)-7-(N-Boc-5-amino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester yielded 27mgs  
 5 (88%) of the desired product, purified by HPLC (Column: Dynamax 60A C18, Method: 10-50% MeCN/H<sub>2</sub>O 10ml/min. Detector: 254nm. Time(%MeCN): 0:00(10), 00:15(10, inj., detect on), 10:00(10), 40:00(50), 55:00(50), 59:58(10, detect off), 60:00(off), R<sub>f</sub>=32.1min). <sup>1</sup>H NMR (D<sub>2</sub>O) 7.30 (1H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, ArH C6-H), 7.24 (1H, dd, <sup>4</sup>J<sub>HH</sub>=2Hz, <sup>3</sup>J<sub>HH</sub>=8Hz, ArH C8-H), 7.08 (1H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, ArH C9-H), 3.85 (1.6H, m, C3-H, NCHHCH(Me)CO<sub>2</sub>), 3.57 (1.8H, m, C3-H, NCHHCH(Me)CO<sub>2</sub>), 3.25 (0.6H, dd, <sup>2</sup>J<sub>HH</sub>=14Hz, <sup>3</sup>J<sub>HH</sub>=5Hz), 3.10 (3H, s, NCH<sub>3</sub>),  
 10 2.74 (3H, m, CH(Me)CO<sub>2</sub>, CH<sub>2</sub>NH<sub>2</sub>), 2.40 (2H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, ArCH<sub>2</sub>), 1.38 (4H, m), 1.13 (4H, m), 0.95 (1.8H, d, <sup>3</sup>J<sub>HH</sub>=7Hz, CHCH<sub>3</sub>), 0.88 (1.2H, d, <sup>3</sup>J<sub>HH</sub>=7Hz, CHCH<sub>3</sub>). Exact mass (FAB, M+H<sup>+</sup>)calculated for C<sub>29</sub>H<sub>39</sub>N<sub>3</sub>O<sub>4</sub> 376.2236, found 376.2228.

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## Example 32



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( $\pm$ )-1-methyl-4-(N-Cbz-2-carboxy-2-aminoethyl)-7-(6-amino-1-hexyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate.

a) A solution of ( $\pm$ )-2-(N-Cbz-amino)-3-aminopropionic acid (1gram, 4.22mmol), 10mL  
 25 methanol, 1.6mL of a 4N HCl in dioxane was stirred at room temperature for 24 hrs. The reaction mixture was concentrated *in vacuo*, triturated with ether (50mL). The resulting residue was dissolved in dimethylformamide (6mL) and allowed to react with N-methyl-6-iodoisatoic anhydride (1.34grams, 4.5mmol) and triethylamine (1.9mL, 14mmol). The reaction was heated to 50°C for 1.5hrs, cooled to room temperature, diluted with 100mL water and extracted with 2X75mL ethyl acetate. The combined organics were dried over sodium  
 30 sulphate, decanted, and concentrated *in vacuo*. The resulting residue was purified by column chromatography (SiO<sub>2</sub>, 35%ethyl acetate/hexane to 65%ethyl acetate/hexane) to yield ( $\pm$ )-N<sup>2</sup>-(N<sup>3</sup>-methyl-2-amino-5-iodobenzoyl)-3-amino-2-(N<sup>1</sup>-Cbz)-aminopropionate methyl ester (1.65 grams, 77%).

- b) ( $\pm$ )-1-methyl-4-(2-carboxy-2-(N-Cbz)-aminoethyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester was prepared using the method described in part (d) of example 31. Thus, 0.825 grams (1.6mmol) of ( $\pm$ )-N<sup>2</sup>-(N<sup>3</sup>-methyl-2-amino-5-iodobenzoyl)-3-amino-2-(N<sup>1</sup>-Cbz)-aminopropionate methyl ester yielded 0.756 grams (86%) of ( $\pm$ )-1-methyl-4-(2-carboxy-2-(N-Cbz)-aminoethyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester (TLC, SiO<sub>2</sub>, 1:1 ethyl acetate/hexane, R<sub>f</sub>=0.21, uv positive). <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 8.09 (0.5H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, C6-H), 8.04 (0.5H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, C6-H), 7.76 (1H, dd, <sup>4</sup>J<sub>HH</sub>=2Hz, <sup>3</sup>J<sub>HH</sub>=8Hz, C8-H), 7.25 (5H, m, Ph), 6.92 (0.5H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, C9-H), 6.87 (0.5H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, C9-H), 6.22 (0.5H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, CbzNH), 5.90 (0.5H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, CbzNH), 4.67 (1H, bs, NHBoc), 5.05 (2H, OCH<sub>2</sub>), 4.63 (1H, m, CHCO<sub>2</sub>), 4.2-3.6 (7H, m, C3-H, NCH<sub>2</sub>CH, C3-H, OCH<sub>3</sub>), 3.29 (1.5H, bs, NCH<sub>3</sub>), 3.24 (1.5H, bs, NCH<sub>3</sub>); <sup>13</sup>C NMR 170.5, 170.4, 168.4, 168.3, 166.8, 166.7, 156.1, 141.3, 140.7, 139.6, 139.4, 136.5, 136.2, 129.8, 128.7, 128.6, 128.5, 128.4, 128.3, 128.1, 123.1, 122.9, 89.5, 76.9, 67.0, 52.6, 52.4, 35.0.
- c) ( $\pm$ )-1-methyl-4-(2-carboxy-2-(N-Cbz)aminoethyl)-7-(N-Boc-5-amino-1-pentynyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester was prepared using the method described in part (f) of example 1. Thus, 0.2 grams (0.36mmol) of ( $\pm$ )-1-methyl-4-(2-carboxy-2-(N-Cbz)-aminoethyl)-7-iodo-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester yielded 0.187 grams (85%) of the desired product (TLC, 1:1 ethyl acetate/hexane R<sub>f</sub>=0.17 uv and ninhydrin positive). <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 7.83 (0.5H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, C6-H), 7.79 (0.5H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, C6-H), 7.50 (1H, bd, <sup>3</sup>J<sub>HH</sub>=8Hz, C8-H), 7.30 (5H, m, Ph), 7.10 (0.5H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, C9-H), 7.07 (0.5H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, C9-H), 6.24 (0.5H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, CbzNH), 5.84 (0.5H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, CbzNH), 5.09 (2H, OCH<sub>2</sub>), 4.66 (2H, m, NHBoc, CHCO<sub>2</sub>), 4.2-3.6 (7H, m, C3-H, NCH<sub>2</sub>CH, C3-H, OCH<sub>3</sub>), 3.34 (1.5H, bs, NCH<sub>3</sub>), 3.29 (1.5H, bs, NCH<sub>3</sub>) 3.24 (2H, m, CH<sub>2</sub>NHBoc), 2.43 (2H, m, CCCH<sub>2</sub>), 1.78 (2H, m, CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>), 1.44 (9H, s, C(CH<sub>3</sub>)<sub>3</sub>); <sup>13</sup>C NMR 170.3, 170.2, 168.2, 168.1, 167.5, 167.4, 155.9, 139.7, 135.1, 135.0, 133.9, 133.8, 128.4, 123.3, 128.0, 127.9, 127.8, 121.4, 120.9, 90.9, 79.1, 79.0, 66.9, 66.8, 53.7, 52.9, 52.8, 52.7, 51.9, 49.6, 34.8, 28.7, 28.3, 16.8.
- d) ( $\pm$ )-1-methyl-4-(2-carboxy-2-(N-Cbz)aminoethyl)-7-(N-Boc-5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester was prepared using the method described in part (a) of example 3. Thus, 93mgs of ( $\pm$ )-1-methyl-4-(2-carboxy-2-(N-Cbz)aminoethyl)-7-(N-Boc-5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester yielded 93mgs (99%) of the desired product. <sup>1</sup>H NMR (CDCl<sub>3</sub>, dTMS) 7.60 (0.6H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, C6-H), 7.55 (0.4H, d, <sup>4</sup>J<sub>HH</sub>=2Hz, C6-H), 7.30 (6H, m, Ph, C8-H), 7.07 (0.6H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, C9-H), 7.06 (0.4H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, C9-H), 6.34 (0.6H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, CbzNH), 5.89 (0.4H, d, <sup>3</sup>J<sub>HH</sub>=8Hz, CbzNH), 5.09 (2H, m, OCH<sub>2</sub>), 4.60 (2H, bs, CHCO<sub>2</sub>, NHBoc), 4.2-3.6 (7H, m, C3-H, NCH<sub>2</sub>CH, C3-H, OCH<sub>3</sub>), 3.33 (1.5H, bs, NCH<sub>3</sub>), 3.30 (1.5H,



bs, NCH<sub>3</sub>) 3.07 (2H, m, CH<sub>2</sub>NHBoc), 2.60 (2H, m, ArCH<sub>2</sub>), 1.60 (2H, m), 1.45 (2H, m), 1.40 (9H, s, C(CH<sub>3</sub>)<sub>3</sub>), 1.33 (2H, m); <sup>13</sup>C NMR 170.4, 168.5, 168.4, 155.9, 140.3, 138.7, 132.4, 130.4, 128.4, 128.3, 128.0, 121.0, 66.9, 60.3, 53.9, 53.0, 52.8, 52.7, 52.0, 50.5, 49.7, 40.4, 34.8, 30.7, 29.9, 28.3, 26.3, 20.9, 14.1.

5

e) (+)-1-methyl-4-(2-carboxy-2-(N-Cbz)aminoethyl)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate was prepared using the method described in part (g) of example 1. Thus 46mgs of (±)-1-methyl-4-(2-carboxy-2-(N-Cbz)aminoethyl)-7-(N-Boc-5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione methyl ester yielded 16mgs (40%) of the desired product, purified by HPLC (Column: Dynamax 60A C18, Method: 10-50% MeCN/H<sub>2</sub>O 10ml/min. Detector: 254nm. Time(%MeCN): 0:00(10), 00:15(10,inj.,detect on), 10:00(10), 40:00(50), 55:00(50), 59:58(10, detect off), 60:00(off) R<sub>f</sub>=38.3min). <sup>1</sup>H NMR (D<sub>2</sub>O) 7.38 (2H, bs), 7.18 (3H, m), 6.95 (2H, m), 6.78 (1H, m), 4.94 (2H, m, OCH<sub>2</sub>Ph), 4.0-3.3 (4H, m), 3.11 (1H, s, NCH<sub>3</sub>), 2.86 (2H, s, NCH<sub>3</sub>), 2.81 (2H, m, CH<sub>2</sub>NH<sub>2</sub>), 2.58 (1.3H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, ArCH<sub>2</sub>), 2.47 (0.7H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, ArCH<sub>2</sub>), 1.57 (4H, m), 1.23 (2H, m).

f) (+)-1-methyl-4-(2-carboxy-2-(N-Cbz)aminoethyl)-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate was prepared using the method described in example 2. Thus 8mgs of (±)-1-methyl-4-(2-carboxy-2-(N-Cbz)aminoethyl)-7-(5-amino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate yielded 6mgs (69%) of (±)-1-methyl-4-(2-carboxy-2-(N-Cbz)aminoethyl)-7-(5-guanidino-1-pentyl)-3,4-dihydro-1H-1,4-benzodiazepine-2,5-dione trifluoroacetate (HPLC, Column: Dynamax 60A C18, Method: 10-50% MeCN/H<sub>2</sub>O 10ml/min. Detector: 254nm. Time(%MeCN): 0:00(10), 00:15(10,inj.,detect on), 10:00(10), 40:00(50), 55:00(50), 59:58(10, detect off), 60:00(off), R<sub>f</sub>=45.4min). <sup>1</sup>H NMR (D<sub>2</sub>O) 7.40 (2H, bs), 7.20 (3H, m), 6.96 (2H, m), 6.78 (1H, t, <sup>3</sup>J<sub>HH</sub>=8Hz), 4.98 (2H, m, OCH<sub>2</sub>Ph), 4.0-3.3 (4H, m), 3.13 (1H, s, NCH<sub>3</sub>), 2.94 (2H, m, CH<sub>2</sub>NH), 2.83 (2H, s, NCH<sub>3</sub>), 2.59 (1.3H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, ArCH<sub>2</sub>), 2.48 (0.7H, t, <sup>3</sup>J<sub>HH</sub>=7Hz, ArCH<sub>2</sub>), 1.42 (4H, m), 1.20 (2H, m). Exact mass calculated for C<sub>27</sub>H<sub>35</sub>N<sub>6</sub>O<sub>6</sub> 539.2618, found 539.2618.

### Example 33

#### Fibrinogen-GP IIb/III<sub>a</sub> Receptor ELISA Binding Assay

35 The method used is essentially that described in Nachman and Leung (J. Clin. Invest., 69: 263-269 [1982]). The GP IIb/III<sub>a</sub> is essentially purified as described in Fitzgerald *et al.*, (Anal. Biochem. 151: 169-177 [1985]).

A. GP IIb/IIIa Purification

Outdated human platelets are washed 3 times with 10 mM tris-HCl, 150 mM NaCl (TBS), 1 mM EDTA, pH 7.5, and centrifuged at 2000 x g to pellet cells. Cells are lysed in 5 pellet volumes of TBS, 1% Triton X-100, 1 mM Ca<sub>2</sub>Cl<sub>2</sub>, and followed by centrifugation at 5 30,000 x g. The supernatant fraction is collected and the supernatant is loaded onto a concanavalin-A column, previously equilibrated in TBS, 1 mM Ca<sub>2</sub>Cl<sub>2</sub>, 0.1% Triton, 0.05% NaN<sub>3</sub> and eluted with 0.2 M  $\alpha$ -methylmannoside. Fractions are pooled and loaded onto a heparin-agarose column. The flowthrough is collected and concentrated on an Amicon YM 30 filter to a volume of approximately 5-10 ml. The concentrate is then applied to an S-300 10 column (500 ml) and 6 ml fractions are collected. The GP IIb/IIIa containing fractions are collected, pooled, and stored at -80° C.

B. Purification of Low Solubility Fraction of Fibrinogen

The purification of fibrinogen is conducted essentially as described by Lipinska *et al.*, (*J. Lab. Clin. Med.* 507, [1974]). Briefly, a 0.3 % w/v solution of human fibrinogen (Kabi 15 #5302) is dissolved in 150 mM NaCl. Saturated (NH<sub>4</sub>)<sub>2</sub>SO<sub>4</sub> is added dropwise with stirring to the fibrinogen solution to obtain about 16% saturation. The precipitate is spun down in appropriate size bottles at 2000 x g. The supernatant is decanted and the precipitate resuspended in 150 mM NaCl (approximately 50% of the original volume). NH<sub>4</sub>SO<sub>4</sub> is again added dropwise to obtain 16% saturation. The suspension is spun down and the 20 precipitate is resuspended in Tris-saline in a minimal volume (approximately 5% of the original volume). Any remaining insoluble material is spun down at 2000 rpm in a Sorval type centrifuge and the fibrinogen supernatant is decanted and dialyzed overnight at 4°C against Tris-saline. Characterization of the fibrinogen is by the Bradford protein assay, SDS-PAGE, and/or Western blotting using well known standard procedures.

25 C. ELISA Assay

Briefly, 96 well plates are coated (Type Nunc 1 Maxisorp™) with 10 mg/ml purified fibrinogen (100  $\mu$ l/well), and allowed to stand overnight at 4°C. The plates are washed three times with PBS Tween (0.137 M NaCl, 0.003 M KCl, 0.008 M Na<sub>2</sub>HPO<sub>4</sub>, 0.001 M KH<sub>2</sub>PO<sub>4</sub>, pH 7.4 at room temperature, 0.05% Tween-20) and blocked for 1 to 2 hours at room 30 temperature with 200  $\mu$ l/well TNCNT (which is 0.5% BSA, 20mM Tris, pH 7.5 at room temperature, 120mM NaCl, 0.2% NaN<sub>3</sub>, 2mM CaCl<sub>2</sub>, 0.05% Tween 20, 0.5% BSA [Calbiochem RIA grade or better]) on a plate shaker. The plates are again washed three times with PBS/Tween and then 50  $\mu$ l of sample in TNCNT is added. The mixture is incubated for 15 minutes at room temperature on a plate shaker. The stock solution of purified GP IIb/IIIa 35 receptor from human platelets, (0.4 - 1.0 mg/ml GP IIb/IIIa in 0.1% Triton X-100, 1mM CaCl<sub>2</sub>, 20 mM Tris, 150 mM NaCl, 0.05% NaN<sub>3</sub> in 0.3 M N-acetyl glucosamine pH 7.5, stored at -70°C), is reconstituted to about 40 mg/ml in TNCNT. Fifty  $\mu$ l of this diluted GP IIb/IIIa is then added to each well and incubated on a plate shaker at room temperature. After one hour, the plates are washed four times with PBS/Tween and 100  $\mu$ l of a polyclonal or

monoclonal antibody specific for GP IIIa such as AP3 (1mg/ml) (See e.g. Newman *et al.*, *Blood*, 65: 227-232 [1985]) in ELISA buffer (PBS, 0.5% BSA, 0.05% Tween 20, 0.01% Thimerosal) is added. After a one hour incubation at room temperature on a plate shaker, the samples are washed 4 times with PBS/Tween. One hundred ml of GAM-HRP (horse  
5 radish peroxidase conjugate of goat anti-mouse IgG [Pel-Freeze Cat. 715305-1] dissolved in ELISA buffer) previously diluted to 1:10,000 is then added and the samples are incubated 1 hour at room temperature on a plate shaker. The samples are then washed 4 times with PBS/Tween and 100 ml OPD/H<sub>2</sub>O<sub>2</sub> substrate is added (OPD/H<sub>2</sub>O<sub>2</sub> substrate: 10 mg o-phenylenediamine in 15 ml phosphate/citrate buffer, at room temperature and covered with  
10 foil; just before use, 6.25 ml of 30% H<sub>2</sub>O<sub>2</sub> is added to give a final solution of 0.67 mg OPD/ml in 0.0125% H<sub>2</sub>O<sub>2</sub>). (The phosphate/citrate buffer consists of 16 mM Citric Acid, 50 mM Na<sub>2</sub>HPO<sub>4</sub>, pH 5.0). The color develops within about 3 to 20 minutes and the reaction is stopped with 100 ml 1 M H<sub>2</sub>SO<sub>4</sub>. The optical density at 492 nm vs 405 nm is recorded and IC<sub>50</sub> values are determined.

15

#### Example 34

#### Human Vitronectin-Vitronectin Receptor ( $\alpha_v\beta_3$ ) ELISA Assay

##### A. Human Vitronectin Purification

20 Human vitronectin (Vn) is isolated from human plasma and purified by affinity chromatography by the method of Yatohgo *et al.*, (*Cell Structure and Function* 13: 281-292 [1988]).

##### B. Human Vitronectin receptor ( $\alpha_v\beta_3$ ) Purification

25 Human vitronectin receptor (VnR) is purified from human placenta by the method of Pytela *et al.*, (*Methods Enzymol.*, 144: 475 [1987]). Alternatively the  $\alpha_v\beta_3$  receptor can be purified from some cell lines (e.g., human embryonic kidney 293 cells) transfected with DNA sequences for both the  $\alpha_v$  and  $\beta_3$  subunits. The subunits are purified by employing octylglucoside extraction followed by Con-A, Heparin-Sepharose, and S-300  
30 Chromatography.

##### C. Monoclonal Antibodies

Anti-GP IIb/IIIa monoclonal antibodies specific for human GP IIIa are prepared by the method of Newman *et al.* (*Blood*, 65: 227-232 [1985]), or a similar procedure. This mouse Mab  
35 is specific for the  $\beta_3$  subunit of the vitronectin receptor.

Rabbit Fab 2 anti-mouse Fc fragment horse radish peroxidase conjugate (anti-MuFc HRP) is obtained from PelFreeze (cat. no. 715305-1).

**D. ELISA Assay**

Maxisorp microtiter plates are coated with 2 mg/ml human vitronectin dissolved in PBS (50 ml/well) and stored overnight at 4°C. The plates are washed two times with PBS-0.05% Tween-20 (wash buffer) and blocked by incubating with about 150 ml/well of assay buffer (1%, BSA [RIA grade or better] in 50 mM Tris-HCl, 100 mM NaCl, 1mM MgCl<sub>2</sub>, CaCl<sub>2</sub>, MnCl<sub>2</sub> pH 7.4) for 60 minutes. Dilutions of standards are prepared and putative inhibitors (Table 3) are dissolved in assay buffer. The blocked plates are emptied and 25 ml/well of inhibitor or standard solution is added to each well. Twenty-five ml of a 30 mg/ml solution of purified  $\alpha_v\beta_3$  in assay buffer is pipetted into the coated plate. The final concentration of receptor in the assay well is about 15 mg/ml. The plate is incubated on a shaker for 60 minutes. Meanwhile, for each microtiter plate, 6 ml buffer solution containing 1.5 mg/ml of mouse-monoclonal antibody specific for  $\beta_3$  is prepared. To this solution is added ml of the secondary antibody, which is anti-mouse-Fc-HRP antibody conjugate. For example, for one plate, prepare 6 ml of a 1.5 mg/ml mouse Mab solution to which is added 1 ml of anti-mouse-Fc-HRP antibody stock, (this represents a 1:6000 dilution of the antibody - HRP conjugate). This mixture is allowed to incubate during the receptor-inhibitor incubation. The assay plates are washed 4 times with PBS-Tween and 50 ml/well of the antibody mixture is then pipetted into the plate for a 60 minute incubation. The plate is washed 4 times and the color reaction is developed with 50 ml/well of 0.67 mg/ml o-phenyldiamine in PBS containing 0.012% H<sub>2</sub>O<sub>2</sub>. Alternatively, 16 mM citric acid, 50 mM Na<sub>2</sub>PO<sub>4</sub> at pH 5.0 can be used as a substrate buffer. The reaction is stopped with 50 ml/well 1 M H<sub>2</sub>SO<sub>4</sub>. The plates are read at 492-405 nm and the data analyzed by four-parameter fit.

**Example 35****GP IIb/IIIa-von Willebrand factor (vWF) ELISA Assay****A. ELISA Assay**

Microtiter plates are coated with 1.0 mg/ml GP IIb/IIIa, prepared by the method of Fitzgerald *et al.*, (*Anal. Biochem.* 151: 169-177 [1985]) and allowed to incubate overnight in coat buffer. The plates are then washed three times in wash buffer (0.05% Tween 20 in PBS) and 150 ml of assay buffer is added and allowed to incubate for 1-2 hours at room temperature on plate shaker. The plates are washed three times and 50 ml of 2x inhibitor in assay buffer (Assay buffer: 0.5% BSA/50mM Tris, 100mM NaCl, 1.0mM CaCl<sub>2</sub>, 1.0mM MgCl<sub>2</sub>, 1.0mM MnCl<sub>2</sub>; coat buffer is the same but without BSA) is added. Fifty ml of 4.0 mg/ml vWF (prepared as described by Ledford *et al.*, *Thrombosis and Haemostasis*, 64(4): 569-575 [1990]) in assay buffer is then added and allowed to incubate for one hour at room temperature on plate-shaker. The plates are washed three times and the antibody mixture is added (1:5000 of mouse anti-vWF and 1:5000 of rabbit-anti-mouse-Fc-HRP, both commercially available) in assay buffer and incubated for 1 hour at room temperature on plate-shaker. Plates are again washed three times and 100 ml of substrate solution (10 mg OPD, 6.5 ml H<sub>2</sub>O<sub>2</sub>, 15 ml

phosphate citrate buffer) is added and incubated at room temperature. The color change of OPD/H<sub>2</sub>O<sub>2</sub> reagent is read at 492 nm with a 405 nm reference wavelength on the filter photometer.

5

## Example 36

## In Vitro Human Platelet Aggregation Assay

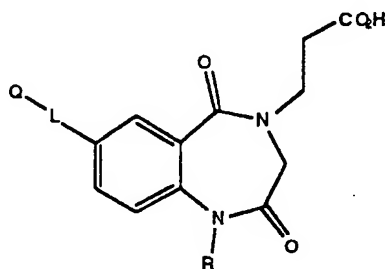
Platelet aggregation assays are performed in human platelet rich plasma (PRP). Fifty milliliters of whole human blood (9 parts) is drawn on 3.6% sodium citrate (1 part) from a donor who has not taken aspirin or related medications for at least two weeks. The blood is centrifuged at 160 x g for 10 minutes at 22°C and allowed to stand for 5 minutes after which the PRP is decanted. Platelet poor plasma (PPP) is isolated from the remaining blood after centrifugation at 2000 x g for 25 minutes. The platelet count of the PRP is diluted to about 300,000 platelets per microliter with PPP.

A 225 ml aliquot of PRP plus 25 ml of either a dilution of the test inhibitor sample or a control (PBS) is incubated for 5 minutes in a Chrono-log Whole Blood Aggregometer at 25°C. An aggregating agent (collagen, 1mg/ml; U46619, 100 ng/ml; or ADP, 17 mM) is added and the transmission is recorded.

20

## Example 37

## Representative 1, 3, 4, and 7 substituted Benzodiazepinediones



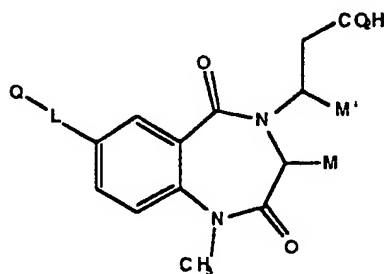
25

Cmpd#	O	L	R
1	amine	(CH <sub>2</sub> ) <sub>3</sub>	H
2	guanidine	(CH <sub>2</sub> ) <sub>5</sub>	H
3	guanidine	(CH <sub>2</sub> ) <sub>3</sub> CC	H
4	guanidine	(CH <sub>2</sub> ) <sub>3</sub>	CH <sub>3</sub>
5	amine	CH <sub>2</sub> CC	CH <sub>3</sub>
6	amine	(CH <sub>2</sub> ) <sub>4</sub>	CH <sub>3</sub>
7	amine	(CH <sub>2</sub> ) <sub>2</sub> CC	CH <sub>3</sub>
8	guanidine	(CH <sub>2</sub> ) <sub>5</sub>	CH <sub>3</sub>

30

	9	guanidine	(CH <sub>2</sub> ) <sub>3</sub> CC	CH <sub>3</sub>
	10	amine	(CH <sub>2</sub> ) <sub>6</sub>	CH <sub>3</sub>
	11	guanidine	(CH <sub>2</sub> ) <sub>4</sub> CC	CH <sub>3</sub>
	12	amine	(CH <sub>2</sub> ) <sub>4</sub> C(O)NH	CH <sub>3</sub>
5	13	guanidine	(CH <sub>2</sub> ) <sub>4</sub> O	CH <sub>3</sub>
	14	amine	CH <sub>2</sub> (pC <sub>6</sub> H <sub>4</sub> )CC	CH <sub>3</sub>
	15	guanidine	(CH <sub>2</sub> ) <sub>5</sub>	(CH <sub>2</sub> ) <sub>4</sub> CH <sub>3</sub>
	16	amine	(CH <sub>2</sub> ) <sub>6</sub>	(CH <sub>2</sub> ) <sub>4</sub> CH <sub>3</sub>
10	17	guanidine	(CH <sub>2</sub> ) <sub>5</sub>	CH <sub>2</sub> (p-ClC <sub>6</sub> H <sub>4</sub> )
	18	amine	(CH <sub>2</sub> ) <sub>3</sub> CC	CH <sub>2</sub> (p-ClC <sub>6</sub> H <sub>4</sub> )
	19	amine	(CH <sub>2</sub> ) <sub>6</sub>	CH <sub>2</sub> (p-ClC <sub>6</sub> H <sub>4</sub> )
	20	amine	(CH <sub>2</sub> ) <sub>4</sub> CC	CH <sub>2</sub> (p-ClC <sub>6</sub> H <sub>4</sub> )
	21	guanidine	(CH <sub>2</sub> ) <sub>7</sub>	CH <sub>2</sub> (p-ClC <sub>6</sub> H <sub>4</sub> )
	22	guanidine	CH <sub>2</sub> CC	CH(C <sub>6</sub> H <sub>5</sub> ) <sub>2</sub>
	23	amine	CH <sub>2</sub> CC	CH(C <sub>6</sub> H <sub>5</sub> ) <sub>2</sub>
	24	guanidine	(CH <sub>2</sub> ) <sub>4</sub>	CH(C <sub>6</sub> H <sub>5</sub> ) <sub>2</sub>
15	25	guanidine	(CH <sub>2</sub> ) <sub>2</sub> CC	CH(C <sub>6</sub> H <sub>5</sub> ) <sub>2</sub>
	26	amine	(CH <sub>2</sub> ) <sub>5</sub>	CH(C <sub>6</sub> H <sub>5</sub> ) <sub>2</sub>
	27	amine	(CH <sub>2</sub> ) <sub>3</sub> CC	CH(C <sub>6</sub> H <sub>5</sub> ) <sub>2</sub>
	28	guanidine	(CH <sub>2</sub> ) <sub>3</sub> CC	CH(C <sub>6</sub> H <sub>5</sub> ) <sub>2</sub>
20	29	amine	(CH <sub>2</sub> ) <sub>6</sub>	CH(C <sub>6</sub> H <sub>5</sub> ) <sub>2</sub>
	30	guanidine	(CH <sub>2</sub> ) <sub>5</sub>	CH <sub>2</sub> (b-naph)
	31	amine	(CH <sub>2</sub> ) <sub>6</sub>	CH <sub>2</sub> (b-naph)
	32	amine	(CH <sub>2</sub> ) <sub>5</sub>	CH <sub>2</sub> (a-naph)
25	33	amine	(CH <sub>2</sub> ) <sub>6</sub>	CH <sub>2</sub> (a-naph)
	34	guanidine	(CH <sub>2</sub> ) <sub>5</sub>	CH <sub>2</sub> (pOMeC <sub>6</sub> H <sub>4</sub> )
	35	amine	(CH <sub>2</sub> ) <sub>6</sub>	CH <sub>2</sub> (pOMeC <sub>6</sub> H <sub>4</sub> )
	36	guanidine	(CH <sub>2</sub> ) <sub>6</sub>	CH <sub>2</sub> (pOMeC <sub>6</sub> H <sub>4</sub> )
	37	amine	(CH <sub>2</sub> ) <sub>4</sub> CC	CH <sub>2</sub> (pNO <sub>2</sub> C <sub>6</sub> H <sub>4</sub> )
	38	guanidine	(CH <sub>2</sub> ) <sub>4</sub> CC	CH <sub>2</sub> (pNO <sub>2</sub> C <sub>6</sub> H <sub>4</sub> )
	39	guanidine	(CH <sub>2</sub> ) <sub>3</sub> CC	CH <sub>2</sub> (mNO <sub>2</sub> C <sub>6</sub> H <sub>4</sub> )
	40	amine	(CH <sub>2</sub> ) <sub>4</sub> CC	CH <sub>2</sub> (mNO <sub>2</sub> C <sub>6</sub> H <sub>4</sub> )
30	41	guanidine	(CH <sub>2</sub> ) <sub>4</sub> CC	CH <sub>2</sub> (mNO <sub>2</sub> C <sub>6</sub> H <sub>4</sub> )
	42	guanidine	(CH <sub>2</sub> ) <sub>5</sub>	CH <sub>2</sub> (pCF <sub>3</sub> C <sub>6</sub> H <sub>4</sub> )
	43	amine	(CH <sub>2</sub> ) <sub>6</sub>	CH <sub>2</sub> (pCF <sub>3</sub> C <sub>6</sub> H <sub>4</sub> )
	44	amine	(CH <sub>2</sub> ) <sub>5</sub>	CH <sub>2</sub> (mOPhC <sub>6</sub> H <sub>4</sub> )
35	45	guanidine	(CH <sub>2</sub> ) <sub>5</sub>	CH <sub>2</sub> (mOPhC <sub>6</sub> H <sub>4</sub> )
	46	amine	(CH <sub>2</sub> ) <sub>6</sub>	CH <sub>2</sub> (mOPhC <sub>6</sub> H <sub>4</sub> )
	47	amine	(CH <sub>2</sub> ) <sub>3</sub> CC	CH <sub>2</sub> (o,pF <sub>2</sub> C <sub>6</sub> H <sub>4</sub> )

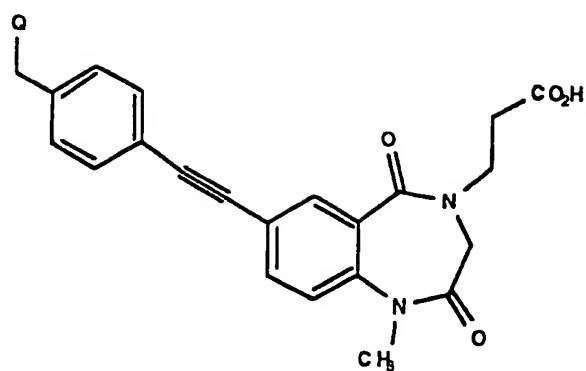
5	48	guanidine	(CH <sub>2</sub> ) <sub>3</sub> CC	CH <sub>2</sub> ( <i>o,p</i> F <sub>2</sub> C <sub>6</sub> H <sub>4</sub> )
	49	amine	(CH <sub>2</sub> ) <sub>4</sub> CC	CH <sub>2</sub> ( <i>o,p</i> F <sub>2</sub> C <sub>6</sub> H <sub>4</sub> )
	50	guanidine	(CH <sub>2</sub> ) <sub>4</sub> CC	CH <sub>2</sub> ( <i>o,p</i> F <sub>2</sub> C <sub>6</sub> H <sub>4</sub> )
	51	amine	(CH <sub>2</sub> ) <sub>5</sub>	CH <sub>2</sub> ( <i>o,p</i> F <sub>2</sub> C <sub>6</sub> H <sub>4</sub> )
	52	guanidine	(CH <sub>2</sub> ) <sub>5</sub>	CH <sub>2</sub> ( <i>o,p</i> F <sub>2</sub> C <sub>6</sub> H <sub>4</sub> )
10	53	amine	(CH <sub>2</sub> ) <sub>6</sub>	CH <sub>2</sub> ( <i>o,p</i> F <sub>2</sub> C <sub>6</sub> H <sub>4</sub> )
	54	guanidine	(CH <sub>2</sub> ) <sub>6</sub>	CH <sub>2</sub> ( <i>o,p</i> F <sub>2</sub> C <sub>6</sub> H <sub>4</sub> )
	55	amine	(CH <sub>2</sub> ) <sub>5</sub>	Ph
	56	guanidine	(CH <sub>2</sub> ) <sub>5</sub>	Ph
	57	amine	(CH <sub>2</sub> ) <sub>6</sub>	Ph
15	58	amine	(CH <sub>2</sub> ) <sub>6</sub>	C <sub>6</sub> H <sub>4</sub> pOMe
	59	amine	(CH <sub>2</sub> ) <sub>5</sub>	cyclohexyl
	60	guanidine	(CH <sub>2</sub> ) <sub>5</sub>	cyclohexyl
	61	amine	(CH <sub>2</sub> ) <sub>6</sub>	cyclohexyl
	62	amine	(CH <sub>2</sub> ) <sub>6</sub>	CH <sub>2</sub> ( <i>p</i> OPhC <sub>6</sub> H <sub>4</sub> )



	Cmpd#	O	L	M	M'
20	63	guanidine	(CH <sub>2</sub> ) <sub>2</sub> CC	CH <sub>3</sub>	H
	64	guanidine	(CH <sub>2</sub> ) <sub>3</sub>	Ph	H
	65	guanidine	CCCH <sub>2</sub>	Ph	H
	66	guanidine	(CH <sub>2</sub> ) <sub>3</sub> CC	H	CH <sub>3</sub>
25	67	amine	(CH <sub>2</sub> ) <sub>6</sub>	H	( <i>R</i> )-CH <sub>3</sub>
	68	amine	(CH <sub>2</sub> ) <sub>6</sub>	H	( <i>R,S</i> )-CH <sub>3</sub>
	69	amine	(CH <sub>2</sub> ) <sub>6</sub>	H	Ph
	70	guanidine	(CH <sub>2</sub> ) <sub>4</sub> CC	H	Ph
	71	amine	(CH <sub>2</sub> ) <sub>6</sub>	H	( <i>R</i> )- <i>p</i> OMeC <sub>6</sub> H <sub>4</sub>
30	72	amine	(CH <sub>2</sub> ) <sub>6</sub>	H	( <i>R,S</i> )- <i>p</i> OMeC <sub>6</sub> H <sub>4</sub>
	73	amine	(CH <sub>2</sub> ) <sub>3</sub>	-(CH <sub>2</sub> ) <sub>2</sub> -	
	74	guanidine	(CH <sub>2</sub> ) <sub>3</sub>	-(CH <sub>2</sub> ) <sub>2</sub> -	



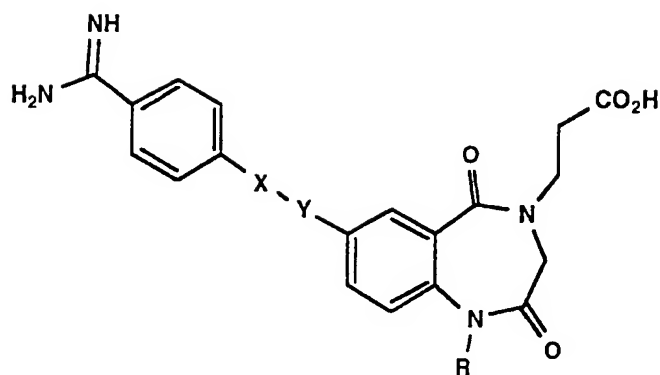




5

Cmpd#	O
84	amine
85	guanidine

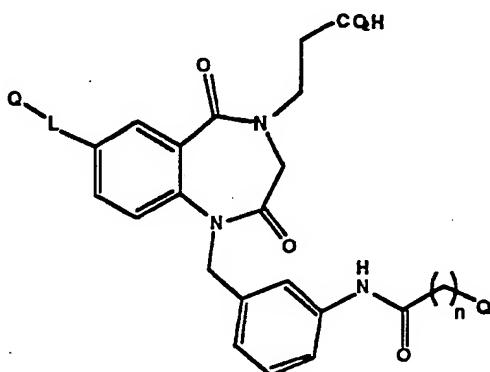
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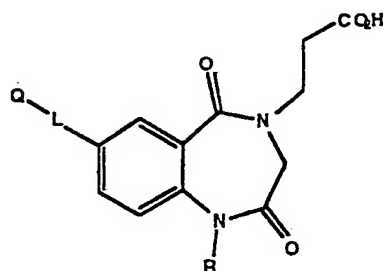
Cmpd#	X	Y	R
86	C	C	CH <sub>3</sub>
87	C	C	CH <sub>2</sub> p-Cl(C <sub>6</sub> H <sub>4</sub> )
88	C(O)	NH	CH <sub>3</sub>
89	CH <sub>2</sub>	O	CH <sub>3</sub>
90	CH <sub>2</sub>	O	CH <sub>2</sub> p-Cl(C <sub>6</sub> H <sub>4</sub> )

20

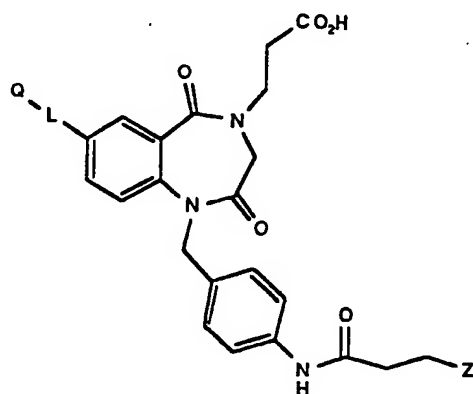


	Cmpd#	O	L	n
5	91	guanidine	(CH <sub>2</sub> ) <sub>3</sub> CC	1
	92	amine	(CH <sub>2</sub> ) <sub>6</sub>	1
	93	amine	(CH <sub>2</sub> ) <sub>6</sub>	2
	94	amine	(CH <sub>2</sub> ) <sub>6</sub>	3
	95	guanidine	(CH <sub>2</sub> ) <sub>6</sub>	3
10	96	amine	(CH <sub>2</sub> ) <sub>6</sub>	4
	97	guanidine	(CH <sub>2</sub> ) <sub>6</sub>	4
	98	amine	(CH <sub>2</sub> ) <sub>6</sub>	5
	99	guanidine	(CH <sub>2</sub> ) <sub>6</sub>	5

15



	Cmpd#	O	L	R
20	100	amine	(CH <sub>2</sub> ) <sub>6</sub>	CH <sub>2</sub> ( <i>m</i> NH <sub>2</sub> C <sub>6</sub> H <sub>4</sub> )
	101	amine	(CH <sub>2</sub> ) <sub>6</sub>	CH <sub>2</sub> ( <i>p</i> NH <sub>2</sub> C <sub>6</sub> H <sub>4</sub> )
	102	amine	(CH <sub>2</sub> ) <sub>6</sub>	CH <sub>2</sub> ( <i>m</i> NHCOBu <sup>n</sup> C <sub>6</sub> H <sub>4</sub> )
	103	amine	(CH <sub>2</sub> ) <sub>6</sub>	(CH <sub>2</sub> ) <sub>5</sub> NH <sub>2</sub>
	104	guanidine	(CH <sub>2</sub> ) <sub>6</sub>	(CH <sub>2</sub> ) <sub>5</sub> NHC(NH)NH <sub>2</sub>
25	105	amine	(CH <sub>2</sub> ) <sub>6</sub>	CH <sub>2</sub> ( <i>m</i> NHCO( <i>pg</i> C <sub>6</sub> H <sub>4</sub> )C <sub>6</sub> H <sub>4</sub> )

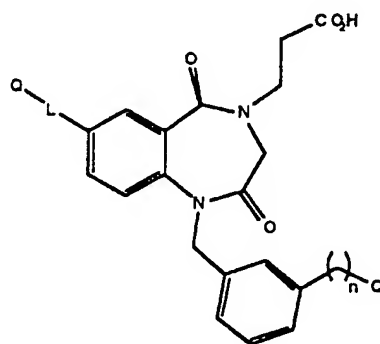


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Cmpd#	Q	L	Z
106	amine	(CH <sub>2</sub> ) <sub>6</sub>	amine
107	guanidine	(CH <sub>2</sub> ) <sub>6</sub>	guanidine
108	amine	(CH <sub>2</sub> ) <sub>6</sub>	Me
109	amine	(CH <sub>2</sub> ) <sub>6</sub>	CO <sub>2</sub> H

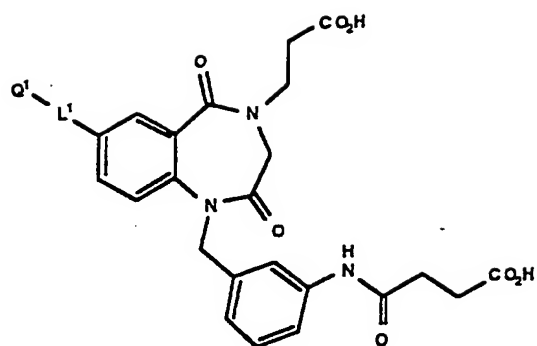
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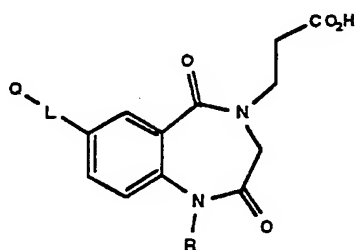


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Cmpd#	Q	L	n
110	amine	(CH <sub>2</sub> ) <sub>6</sub>	4
111	guanidine	(CH <sub>2</sub> ) <sub>6</sub>	4



Compd#	O	L
112	amine	(CH <sub>2</sub> ) <sub>6</sub>
113	guanidine	(CH <sub>2</sub> ) <sub>6</sub>

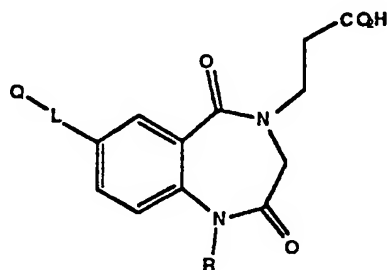


Gmpd#	O	L	R
114	guanidine	(CH <sub>2</sub> ) <sub>5</sub>	(CH <sub>2</sub> ) <sub>4</sub> CO <sub>2</sub> H
115	guanidine	(CH <sub>2</sub> ) <sub>6</sub>	(CH <sub>2</sub> ) <sub>4</sub> CO <sub>2</sub> H
116	guanidine	(CH <sub>2</sub> ) <sub>5</sub>	CH <sub>2</sub> (pCO <sub>2</sub> HC <sub>6</sub> H <sub>4</sub> )
117	amine	(CH <sub>2</sub> ) <sub>6</sub>	CH <sub>2</sub> (pCO <sub>2</sub> HC <sub>6</sub> H <sub>4</sub> )
118	guanidine	(CH <sub>2</sub> ) <sub>6</sub>	CH <sub>2</sub> (pCO <sub>2</sub> HC <sub>6</sub> H <sub>4</sub> )

## Example 38

## Results of ELISA and Platelet Aggregation (PA) Assays for 1, 3, 4, and 7 substituted Benzodiazepinediones

5

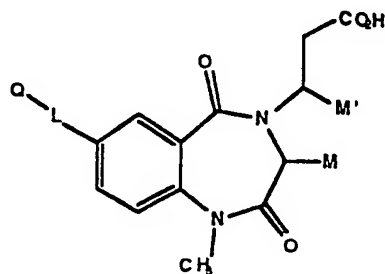


	Cmpd#	IC <sub>50</sub> (mM)	IC <sub>50</sub> (mM)	IC <sub>50</sub> (mM)
		Fg/IIbIIIa	PA(citrate)	PA(heparin)
10	1	8.3		
	2	0.204		
	3	0.254		
	4	0.466		
	5	7.0		
15	6	5.0		
	7	6.0		
	8	0.055	4.58	
	9	0.0036	3.1(2)	7.5
	10	0.070	7.99	
20	11	0.060		
	12	0.030	6.72	
	13	0.008	1.86	
	14	0.012	1.7	3.34
	15	0.011	0.59	
25	16	0.014	1.51	
	17	0.011	2.55	
	18	0.062		
	19	0.014	1.20	4.76
	20	0.014	2.12	
30	21	0.200		
	22	0.010	0.52	
	23	0.056	0.59	
	24	0.013	0.78	1.60
	25	0.015	1.78(2)	

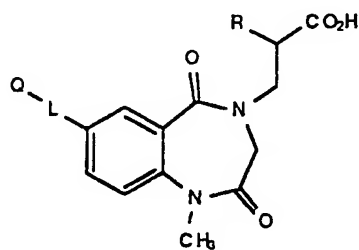
WO 93/08174

PCT/US92/08788

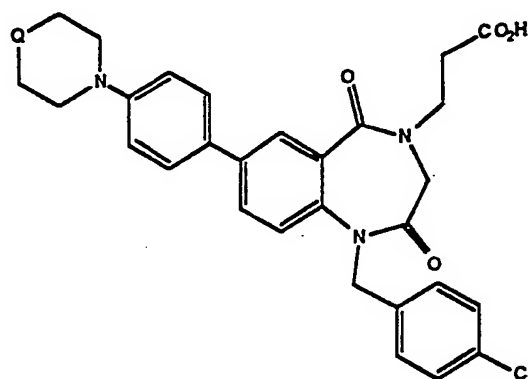
	26	0.020(2)	1.66	3.4
	27	0.0024	0.44	1.3
	28	0.008	0.34	0.95
	29	0.022	0.46	1.1
5	30	0.011(2)	6.88	
	31	0.017	3.19	
	32	0.066	6.53	
	33	0.015	0.76	2.60
	34	0.011	1.12	
10	35	0.023	0.98	2.90
	36	0.024	10.4	
	37	0.031		
	38	0.031		
	39	0.0032	0.94	
15	40	0.008		2.20
	41	0.006		15.0
	42	0.014	2.47	
	43	0.005	1.28	6.30
	44	0.108		
20	45	0.011	18.2	
	46	0.009	18.3	
	47	0.130		
	48	0.024	1.78	
	49	0.027	0.62	
25	50	0.025		
	51	0.200		
	52	0.016	10.0	
	53	0.021	1.73	4.0
	54	0.038		
30	55	0.034		
	56	0.004	0.42	0.99
	57	0.004	0.59	
	58	0.020	0.64	
	59	0.127		
35	60	0.007	1.31	
	61	0.019	1.56	
	62	0.018	17.3	



		IC <sub>50</sub> (mM)	IC <sub>50</sub> (mM)	IC <sub>50</sub> (mM)
	Cmpd#	Fg/IIbIIIa	(citrate)	(heparin)
5	63	80.0		
	64	2.30		
	65	25.0		
	66	0.018	1.19	
	67	0.052(2)	2.3	4.03
10	68	0.040(2)		
	69	0.200		
	70	0.100		
	71	0.097(2)	11.4	41.3
	72	0.190(2)		
15	73	>100.0		
	74	>100.0		

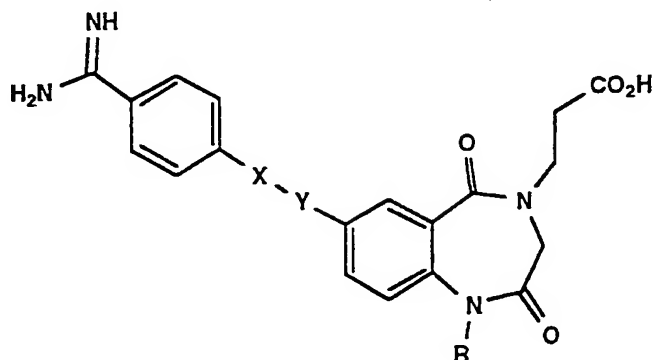


		IC <sub>50</sub> (mM)	IC <sub>50</sub> (mM)	IC <sub>50</sub> (mM)
	Cmpd#	Fg/IIbIIIa	(citrate)	(heparin)
20	75	0.022	2.29	
	76	0.023	2.25	
	77	0.192		
	78	1.27		
25	79	0.149		
	80	0.211		
	81	0.308		



5	Cmpd#	IC <sub>50</sub> (mM)	IC <sub>50</sub> (mM)	IC <sub>50</sub> (mM)
		Fg/IIbIIIa	PA (citrate)	PA (heparin)
	82	0.007	0.67	2.3
	83	0.019	1.80	4.8

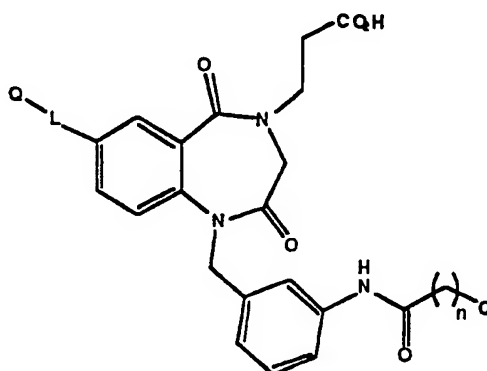
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15	Cmpd#	IC <sub>50</sub> (mM)	IC <sub>50</sub> (mM)	IC <sub>50</sub> (mM)
		Fg/IIbIIIa	PA (citrate)	PA (heparin)
	86	0.035	0.12	0.35
	87	0.007	0.45	0.36
	88	0.008	0.059	0.080
	89	0.005	0.091	0.130
	90	0.015	0.136	

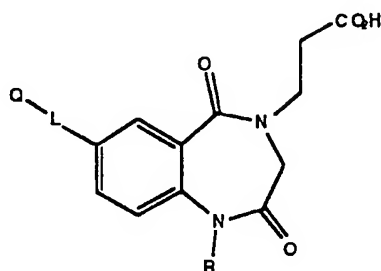
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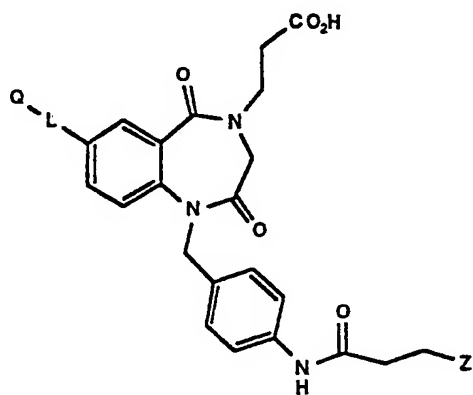


Cmpd#	IC <sub>50</sub> (mM)		
	Fg/IIbIIIa	PA(citrate)	PA(heparin)
5	91	0.018	0.61
	92	0.012	0.55
	93	0.008	0.28
	94	0.008	0.40
	95	0.010	0.87
10	96	0.008	0.31
	97	0.013	0.92
	98	0.011	0.27
	99	0.013	0.64

15

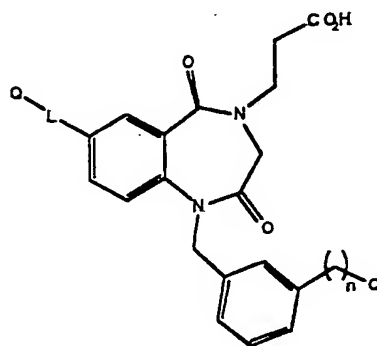


Cmpd#	IC <sub>50</sub> (mM)		
	Fg/IIbIIIa	PA(citrate)	PA(heparin)
20	100	0.012(2)	0.49
	101	0.024	
	102	0.015(2)	0.64
	103	0.064	0.82
	104	0.040	
25	105	0.008	0.35
			0.54



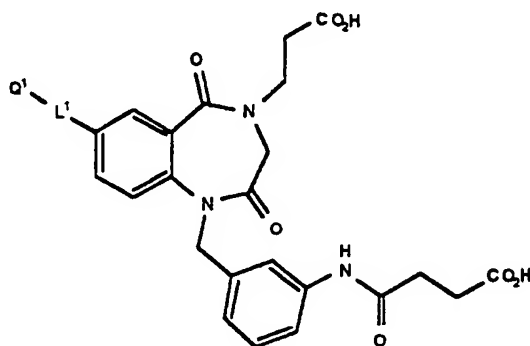
5	Cmpd#	IC <sub>50</sub> (mM)	IC <sub>50</sub> (mM)	IC <sub>50</sub> (mM)
		Fg/IIbIIIa	PA(citrate)	PA(heparin)
	106	0.019	0.13	0.49
	107	0.013		
	108	0.015	0.59	0.89
	109	0.027	1.04	3.4

10



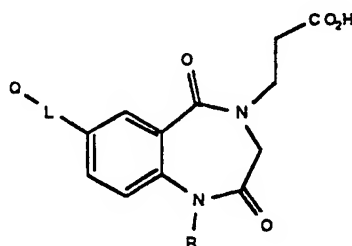
15	Cmpd#	IC <sub>50</sub> (mM)	IC <sub>50</sub> (mM)	IC <sub>50</sub> (mM)
		Fg/IIbIIIa	PA(citrate)	PA(heparin)
	110	0.024	0.70	
	111	0.020	1.65	

20



Cmpd#	IC <sub>50</sub> (mM)	
	Fg/IIbIIIa	PA(citrate)
5 112	0.006	0.48
113	0.012	

10



Cmpd#	IC <sub>50</sub> (mM)		
	Fg/IIbIIIa	PA(citrate)	PA(heparin)
114	0.014	5.56	
15 115	0.053	18.8	
116	0.016	1.68	
117	0.008	1.71	1.80
118	0.019	2.84	

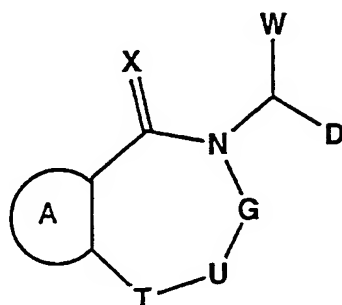
20 While the invention has necessarily been described in conjunction with preferred embodiments, one of ordinary skill, after reading the foregoing specification, will be to effect various changes, substitutions of equivalents, and alterations to the subject matter set forth herein, without departing from the spirit and the scope thereof. Hence, the invention can be practiced in ways other than those specifically described herein. It is therefore intended

25 that the protection granted by Letters Patent hereon be limited only by the appended claims and the equivalents thereof.

What is claimed is:

1. A compound represented by structural formula (I):

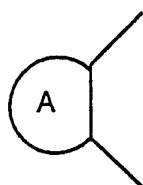
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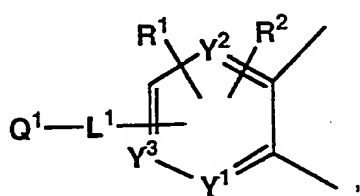
(I)

where the partial structure

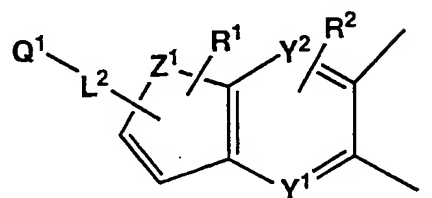
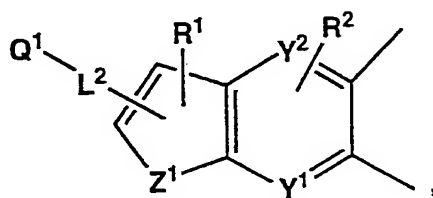
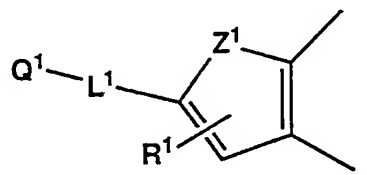
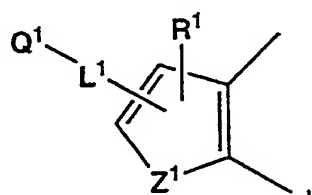
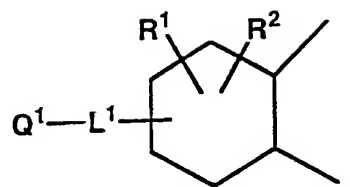
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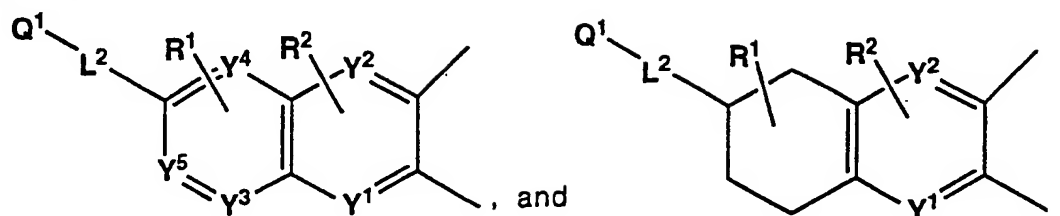
represents



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5 where

$R^1$  and  $R^2$  are independently selected from hydrogen, halo (F, Cl, Br, I), cyano, carboxamido, carbamoyloxy, formyloxy, formyl, azido, nitro, ureido, thioureido, thiocyanato, hydroxy, mercapto, sulfonamido, and an optionally substituted radical selected from C<sub>1</sub>-C<sub>12</sub> alkyl, C<sub>2</sub>-C<sub>12</sub> alkenyl, C<sub>3</sub>-C<sub>12</sub> alkynyl, C<sub>3</sub>-C<sub>12</sub> cycloalkyl, C<sub>6</sub>-C<sub>14</sub> aryl, C<sub>6</sub>-C<sub>10</sub> aryl-C<sub>1</sub>-C<sub>8</sub>-alkyl, C<sub>1</sub>-C<sub>12</sub> alkyloxy, C<sub>6</sub>-C<sub>14</sub> aryloxy, C<sub>1</sub>-C<sub>12</sub> acylamino, N,N-di(C<sub>1</sub>-C<sub>12</sub>)acylamino, N-(C<sub>1</sub>-C<sub>12</sub>)alkyl-N-(C<sub>1</sub>-sulfonylamino, C<sub>1</sub>-C<sub>12</sub> alkylthiocarbonyl, C<sub>1</sub>-C<sub>12</sub> alkylthio, C<sub>1</sub>-C<sub>12</sub> alkylsulfinyl, C<sub>1</sub>-C<sub>12</sub> alkylsulfonyl, C<sub>1</sub>-C<sub>12</sub> alkylsulfonato, N-(C<sub>1</sub>-C<sub>12</sub>)alkylsulfonamido, N,N-di-(C<sub>1</sub>-C<sub>12</sub>) sulfonamido, N-(C<sub>1</sub>-C<sub>12</sub>) alkyl-N-thioformylamino, C<sub>1</sub>-C<sub>12</sub> thioacylamino, N-(C<sub>1</sub>-C<sub>12</sub>)alkyl-N-(C<sub>1</sub>-C<sub>12</sub>) thioacylamino, C<sub>1</sub>-C<sub>12</sub> alkylsulfinamido, N-(C<sub>1</sub>-C<sub>12</sub>)alkyl-N-(C<sub>1</sub>-C<sub>12</sub>)alkylsulfinylamino, C<sub>1</sub>-C<sub>12</sub> carbalkoxy, C<sub>1</sub>-C<sub>12</sub> alkylcarbonyl, C<sub>1</sub>-C<sub>12</sub> alkanoyloxy, N-(C<sub>1</sub>-C<sub>12</sub>)alkylcarboxamido, N,N-di-(C<sub>1</sub>-C<sub>12</sub>)carboxamido, N-(C<sub>1</sub>-C<sub>12</sub>) alkylcarbamoyloxy, N,N-di-(C<sub>1</sub>-C<sub>12</sub>)carbamoyloxy, and heterocycloalkyl or heteroaryl having from 1 to 3 rings, each ring having from 5 to 7 atoms with from 0-3 heteroatoms selected from N, O, and S, provided that at least one ring contains a heteroatom, where the substituents are selected from halo (F, Cl, Br, I), cyano, azido, nitro, hydroxy, mercapto, sulfonamido, ureido, thioureido, carboxamido, carbamoyloxy, formyloxy, formyl, C<sub>1</sub>-C<sub>4</sub>alkyl, C<sub>1</sub>-C<sub>4</sub>alkoxy, phenyl, and phenoxy, optionally,  $R^1$  and  $R^2$  when bonded to adjacent carbon atoms may join to form an unsubstituted or substituted aryl ring, where the substituents are selected from halo (F, Cl, Br, I), cyano, azido, nitro, hydroxy, mercapto, sulfonamido, ureido, thioureido, carboxamido, carbamoyloxy, formyloxy, formyl, C<sub>1</sub>-C<sub>4</sub>alkyl, C<sub>1</sub>-C<sub>4</sub>alkoxy, phenyl, and phenoxy;

30  $Y^1, Y^2, Y^3, Y^4$  and  $Y^5$  are independently selected from CH,  $CR^1$ ,  $CR^2$ , and N;

X is selected from O, S, =NCN, and =NO(C<sub>1</sub>-C<sub>3</sub>)alkyl;

$Z^1$  is selected from CH<sub>2</sub>, NH, S, and O;

35  $Q^1$  is selected from the group consisting of

(A) an amino group selected from

-NH<sub>2</sub>,

-NR<sup>3</sup>H,

40 -NR<sup>3</sup>R<sup>4</sup>, and

-NR<sup>3</sup>R<sup>4</sup>R<sup>5</sup>,

where  $R^3, R^4$ , and  $R^5$  are independently selected from

(i) cyano,

(ii) an optionally substituted radical selected from

45 (a) -NR<sup>6</sup>R<sup>7</sup>,

(b) -C(=NR<sup>8</sup>)-NR<sup>6</sup>R<sup>7</sup>,

(c) -N=CR<sup>9</sup>-NR<sup>6</sup>R<sup>7</sup>,

(d) -NR<sup>10</sup>-CR<sup>9</sup>=NR<sup>8</sup>, and

(e)  $-NR^{10}-C(=NR^8)-NR^6R^7$ ,

where each  $R^6$ ,  $R^7$ ,  $R^8$ ,  $R^9$ , and  $R^{10}$  is independently selected from

hydrogen,

C<sub>1</sub>-C<sub>4</sub> alkoxy,

C<sub>1</sub>-C<sub>4</sub> alkyl, and

halo(F, Cl, Br, I)-C<sub>1</sub>-C<sub>4</sub>-alkyl,

(iii) optionally substituted C<sub>1</sub>-C<sub>12</sub> alkyl,

(iv) optionally substituted C<sub>3</sub>-C<sub>12</sub> cycloalkyl,

(v) optionally substituted C<sub>6</sub>-C<sub>14</sub> aryl,

(vi) optionally substituted C<sub>1</sub>-C<sub>6</sub> alkyl-C<sub>6</sub>-C<sub>14</sub> aryl,

(vii) optionally substituted C<sub>1</sub>-C<sub>8</sub> alkoxy,

(viii) optionally substituted C<sub>6</sub>-C<sub>14</sub> aryloxy, and

(xi) optionally substituted C<sub>1</sub>-C<sub>8</sub> alkoxy-carbonyl,

where the substituents are one to three  $R^{11}$ , each  $R^{11}$  independently selected from

(a) optionally substituted C<sub>6</sub>-C<sub>12</sub> aryloxy,

(b) optionally substituted C<sub>6</sub>-C<sub>12</sub> arylamino,

(c) optionally substituted C<sub>6</sub>-C<sub>12</sub> aroyl,

(d) optionally substituted C<sub>1</sub>-C<sub>8</sub> alkoxy,

(e) optionally substituted C<sub>1</sub>-C<sub>8</sub> alkoxyalkyl,

(f) optionally substituted C<sub>1</sub>-C<sub>8</sub> alkoxyalkyloxy,

(g) optionally substituted C<sub>1</sub>-C<sub>8</sub> alkoxy-carbonyl,

(h) optionally substituted C<sub>1</sub>-C<sub>8</sub> alkyl-carbonyl,

(i) optionally substituted C<sub>1</sub>-C<sub>8</sub> aralkyl-carbonyl,

(j) optionally substituted C<sub>1</sub>-C<sub>8</sub> alkylthio-carbonyl,

(k) optionally substituted C<sub>6</sub>-C<sub>12</sub> aralkylthio-carbonyl,

(l) optionally substituted C<sub>2</sub>-C<sub>8</sub> alkoxythio-carbonyl,

(m) optionally substituted C<sub>6</sub>-C<sub>12</sub> aryl,

(n) optionally substituted C<sub>1</sub>-C<sub>4</sub> alkanoylamino,

(o) optionally substituted C<sub>1</sub>-C<sub>6</sub> alkoxy-carbonyl-C<sub>0</sub>-C<sub>6</sub>alkylamino,

(p) optionally substituted C<sub>1</sub>-C<sub>8</sub> alkylsulfonylamino,

(q) optionally substituted C<sub>6</sub>-C<sub>10</sub> aralkylsulfonylamino,

(r) optionally substituted C<sub>6</sub>-C<sub>10</sub> aralkyl,

(s) optionally substituted C<sub>6</sub>-C<sub>10</sub> alkaryl,

(t) optionally substituted C<sub>1</sub>-C<sub>8</sub> alkylthio,

(u) optionally substituted C<sub>6</sub>-C<sub>10</sub> aralkylthio,

(v) optionally substituted C<sub>1</sub>-C<sub>8</sub> alkylsulfinyl,

(w) optionally substituted C<sub>6</sub>-C<sub>10</sub> aralkylsulfinyl,

(x) optionally substituted C<sub>1</sub>-C<sub>8</sub> alkylsulfonyl,

(y) optionally substituted C<sub>6</sub>-C<sub>10</sub> aralkylsulfonyl,

(z) optionally substituted C<sub>1</sub>-C<sub>8</sub> alkylaminosulfonyl,

(aa) optionally substituted C<sub>6</sub>-C<sub>10</sub> aralkylsulfonylamino,

(ab) optionally substituted 5- or 6-member heterocycle containing 1-4 hetero atoms selected from N, O, and S, and

(ac) optionally substituted C<sub>2</sub>-C<sub>8</sub> alkenyl,

where the substituents are one to three  $R^{12}$ , each  $R^{12}$  independently selected from

nitro,

amino,

C<sub>1</sub>-C<sub>8</sub> alkylamino,

di-(C<sub>1</sub>-C<sub>8</sub>) alkylamino,

amidino,

aminomethyleneimino,

imino,

- 5 imino-C<sub>1</sub>-C<sub>4</sub> alkyl,  
iminomethyleneamino,  
guanidino,  
C<sub>6</sub>-C<sub>10</sub> arylamino,  
C<sub>1</sub>-C<sub>8</sub> acylamino,  
C<sub>1</sub>-C<sub>4</sub> alkylsulfonamino,  
azido,  
cyano,  
hydroxy,  
10 hydroxy-C<sub>1</sub>-C<sub>8</sub>-alkyl,  
C<sub>1</sub>-C<sub>8</sub>-alkoxy,  
phenyloxy,  
C<sub>1</sub>-C<sub>8</sub> alkanoyloxy,  
C<sub>1</sub>-C<sub>8</sub> alkanoyl,  
15 C<sub>6</sub>-C<sub>12</sub> aroyl,  
benzamido,  
phenyl,  
halo(F, Cl, Br, I),  
halo-C<sub>1</sub>-C<sub>8</sub>-alkyl, and  
20 C<sub>1</sub>-C<sub>8</sub>-alkyl,  
(ad) aminosulfonyl,  
(ae) oxo,  
(af) thio,  
(ag) thiocarbonyl,  
25 (ah) hydroxy,  
(ai) mercapto,  
(aj) formyl,  
(ak) formyloxy,  
(al) carboxy,  
30 (am) amino,  
(an) ureido,  
(ao) amidino,  
(ap) guanidino,  
(aq) aminomethyleneimino,  
35 (ar) imino,  
(as) glycyI,  
(at) phthalimido,  
(au) succinimido,  
(av) morpholino,  
40 (aw) C<sub>3</sub>-C<sub>7</sub> cycloalkyl, and  
(ax) halo(F, Cl, Br, I),  
optionally R<sup>3</sup> and R<sup>4</sup> taken together may form optionally substituted  
tetramethylene,  
pentamethylene,  
45 3-oxopentamethylene, and  
3-azapentamethylene,  
where the substituents are selected from one to three R<sup>12</sup>,  
(B) an amidino group selected from  
-C(=NH)-NH<sub>2</sub>,  
50 -C(=NH)-NHR<sup>3</sup>,  
-C(=NR<sup>4</sup>)-NHR<sup>3</sup>,  
-C(=NH)-NR<sup>3</sup>R<sup>4</sup>, and  
-C(=NR<sup>5</sup>)-NR<sup>3</sup>R<sup>4</sup>,  
where R<sup>3</sup>, R<sup>4</sup>, and R<sup>5</sup> are defined above,  
55 (C) an aminoalkyleneimino group selected from  
-N=CH-NH<sub>2</sub>,

-N=CH-NHR<sup>3</sup>,  
 -N=CH-NR<sup>3</sup>R<sup>4</sup>, and  
 -N=CR<sup>5</sup>-NR<sup>3</sup>R<sup>4</sup>,

where R<sup>3</sup>, R<sup>4</sup>, and R<sup>5</sup> are defined above,  
 (D) an iminoalkyleneamino group selected from

-NH-CH=NH,  
 -NH-CH=NR<sup>3</sup>,  
 -NH-CR<sup>4</sup>=NR<sup>3</sup>, and  
 -NR<sup>5</sup>-CR<sup>4</sup>=NR<sup>3</sup>,

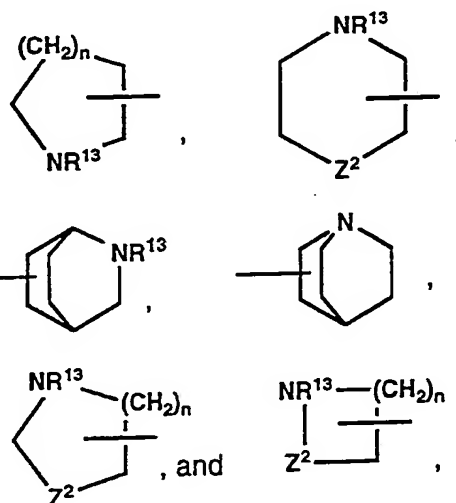
where R<sup>3</sup>, R<sup>4</sup>, and R<sup>5</sup> are defined above,  
 (E) a guanidino group selected from

-NH-C(=NH)-NH<sub>2</sub>,  
 -NH-C(=NH)-NR<sup>3</sup>H,  
 -NH-C(=NH)-NR<sup>3</sup>R<sup>4</sup>,  
 -NH-C(=NR<sup>5</sup>)-NR<sup>3</sup>R<sup>4</sup>,  
 -NR<sup>3</sup>-C(=NR<sup>3</sup>)-NR<sup>3</sup>R<sup>4</sup>,  
 -NR<sup>3</sup>-C(=NH)-NR<sup>3</sup>R<sup>4</sup>,  
 -NR<sup>3</sup>-C(=NR<sup>3</sup>)-NH<sub>2</sub>,  
 -NR<sup>3</sup>-C(=NH)-NH<sub>2</sub>,

-NR<sup>3</sup>-C(=NR<sup>3</sup>)-NHR<sup>4</sup>, and  
 -NR<sup>3</sup>-C(=NH)-NHR<sup>4</sup>,

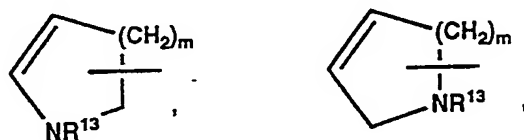
where R<sup>3</sup>, R<sup>4</sup>, and R<sup>5</sup> are defined above;

(F) an optionally substituted saturated heterocyclic group selected from

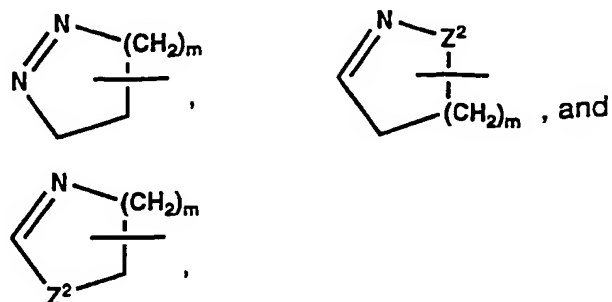


where (1)  $n$  is 0, 1, 2, or 3, (2)  $R^{13}$  is selected from  $R^6$ ,  $-CR^9=NR^8$ ,  $-CR^9(=NR^8)-NR^6R^7$ ,  $-C(=NR^8)-NR^6R^7$ ,  $-N=CR^9-NR^6R^7$ ,  $-NR^{10}-CR^9=NR^8$ , and  $-NR^{10}-C(=NR^8)-NR^6R^7$  where  $R^6$ - $R^{10}$  are defined above, (3)  $Z^2$  is O, S, or  $NR^{13}$ , and (4) the substituents are independently one to three  $R^{12}$ ;

(G) an optionally substituted unsaturated (nonaromatic) heterocyclyl selected from



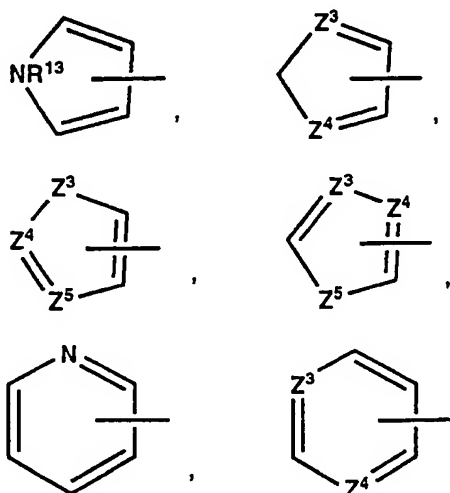




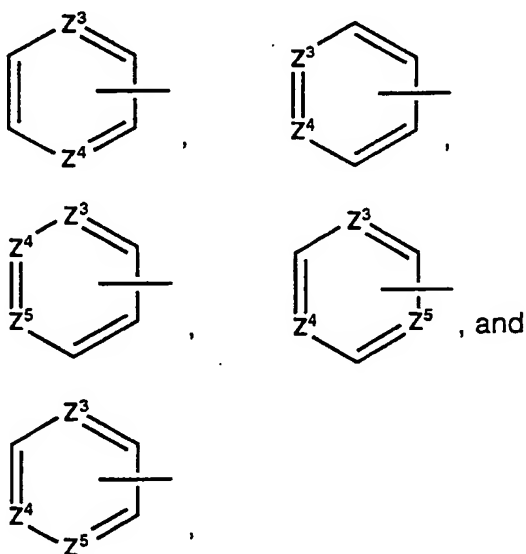
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where (1)  $m$  is 1, 2, or 3, (2)  $\text{Z}^2$  and  $\text{R}^{13}$  are defined above, and (3) the substituents are independently one to three  $\text{R}^{12}$ ;  
 (H) an optionally substituted unsaturated (aromatic) heterocyclyl selected from

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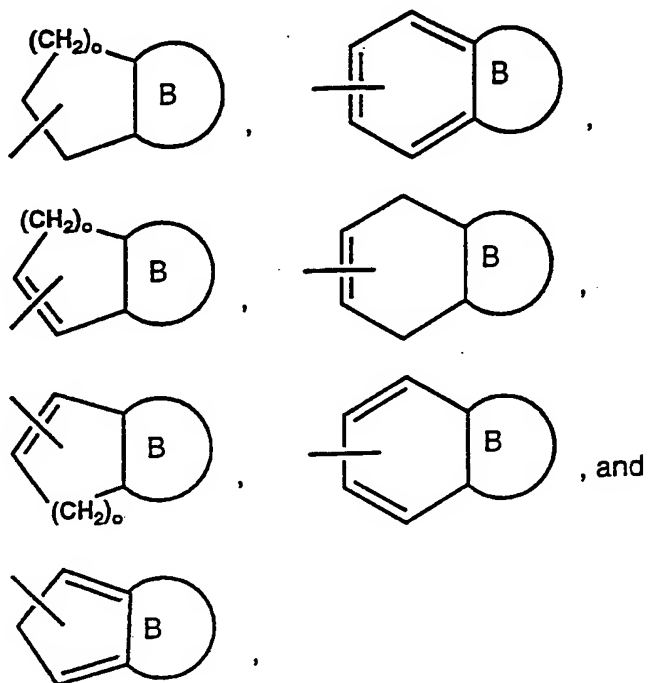
where (1)  $\text{Z}^3$ ,  $\text{Z}^4$ , and  $\text{Z}^5$  are selected from O, S, N, and NH, provided that at least one  $\text{Z}^3$ ,  $\text{Z}^4$ , or  $\text{Z}^5$  is N or NH, (2)  $\text{R}^{13}$  is defined above, and the substituents are independently one to three  $\text{R}^{12}$ ,  
 (I) an optionally substituted bicycloheterocyclic group selected from;

5

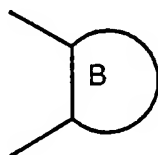
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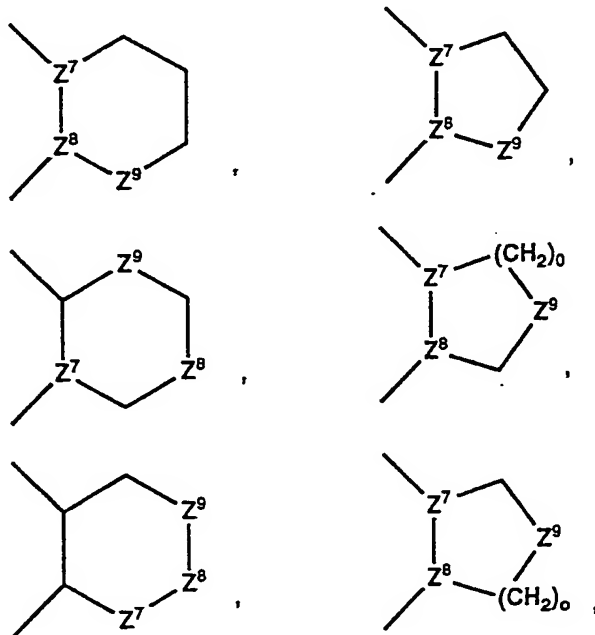
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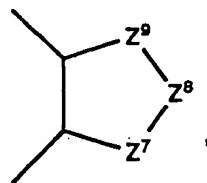
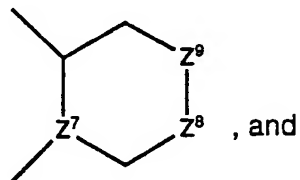
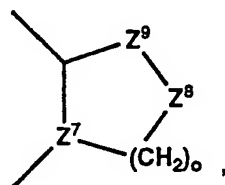
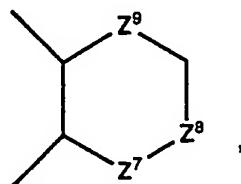
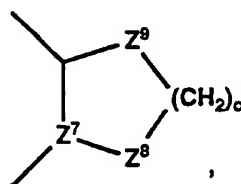
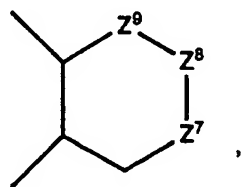
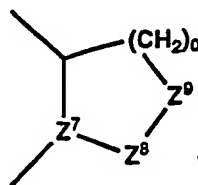
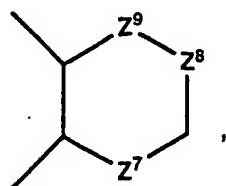
where the partial structure



represents

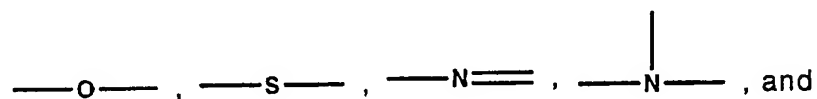
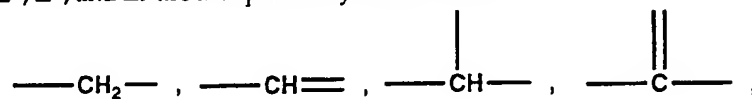


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where  $Z^7$ ,  $Z^8$ , and  $Z^9$  are independently selected from

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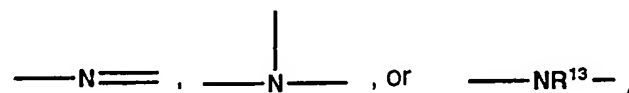


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provided that at least one  $Z^7$ ,  $Z^8$ , or  $Z^9$  is

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where (1)  $o$  is 0, 1, or 2, (2)  $R^{13}$  is defined above, and (3) the substituents are independently one to three  $R^{12}$ ;

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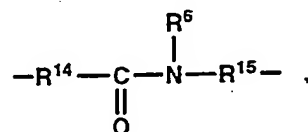
$L^1$  is an optionally substituted bivalent radical selected from the group consisting of

- (A) C<sub>3</sub>-C<sub>7</sub>-alkylene,
- (B) C<sub>3</sub>-C<sub>7</sub>-cycloalkylene,
- (C) C<sub>3</sub>-C<sub>7</sub>-alkenylene,

- (D) C<sub>3</sub>-C<sub>7</sub>-alkadienylene,  
 (E) C<sub>3</sub>-C<sub>7</sub>-alkynylene,  
 (F) C<sub>4</sub>-C<sub>7</sub>-alkadiynylene,  
 (G) C<sub>4</sub>-C<sub>7</sub>-alkenynylene,  
 5 (H) C<sub>6</sub>-C<sub>14</sub>-arylene,  
 (I) C<sub>6</sub>-C<sub>14</sub>-aryl-C<sub>2</sub>-C<sub>4</sub>-alkynylene,  
 (J) C<sub>1</sub>-C<sub>3</sub>-alkyl-C<sub>6</sub>-C<sub>14</sub>-aryl-C<sub>2</sub>-C<sub>4</sub>-alkynylene,  
 (K) C<sub>6</sub>-C<sub>14</sub>-aryl-C<sub>2</sub>-C<sub>4</sub>-alkenylene,  
 (L) C<sub>1</sub>-C<sub>3</sub>-alkyl-C<sub>6</sub>-C<sub>14</sub>-arylene,  
 10 (M) C<sub>1</sub>-C<sub>3</sub>-alkyl-C<sub>6</sub>-C<sub>14</sub>-aryl-C<sub>2</sub>-C<sub>4</sub>-alkenylene,  
 (N) C<sub>6</sub>-C<sub>14</sub>-aryl-C<sub>1</sub>-C<sub>3</sub>-alkylene,  
 (O) C<sub>6</sub>-C<sub>14</sub>-aryl-C<sub>1</sub>-C<sub>3</sub>-alkyloxyene,  
 (P) C<sub>1</sub>-C<sub>2</sub>-alkyl-C<sub>6</sub>-C<sub>14</sub>-aryl-C<sub>1</sub>-C<sub>2</sub>-alkylene,  
 (Q) C<sub>1</sub>-C<sub>3</sub>-alkyloxy-C<sub>6</sub>-C<sub>14</sub>-arylene,  
 15 (R) C<sub>2</sub>-C<sub>6</sub>-alkyloxyene,  
 (S) C<sub>1</sub>-C<sub>5</sub>-alkyloxy-C<sub>1</sub>-C<sub>5</sub>-alkylene,  
 (T) C<sub>6</sub>-C<sub>10</sub>-aryloxyene,  
 (U) C<sub>6</sub>-C<sub>10</sub>-aryloxy-C<sub>1</sub>-C<sub>5</sub>-alkylene,  
 (V) C<sub>2</sub>-C<sub>6</sub>-alkylthioene,  
 20 (W) C<sub>1</sub>-C<sub>5</sub>-alkylthio-C<sub>1</sub>-C<sub>5</sub>-alkylene,  
 (X) C<sub>6</sub>-C<sub>10</sub>-arylthioene,  
 (Y) C<sub>6</sub>-C<sub>10</sub>-arylthio-C<sub>1</sub>-C<sub>5</sub>-alkylene,  
 (Z) C<sub>1</sub>-C<sub>5</sub>-alkylsulfoxide-C<sub>1</sub>-C<sub>5</sub>-alkylene,  
 (AA) C<sub>1</sub>-C<sub>5</sub>-alkylsulfone-C<sub>1</sub>-C<sub>5</sub>-alkylene,

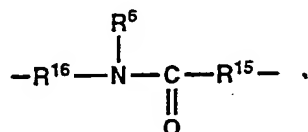
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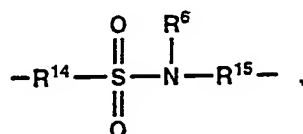


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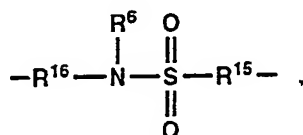


(AD)

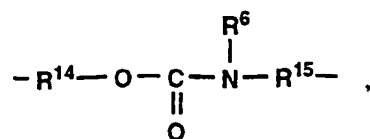


35

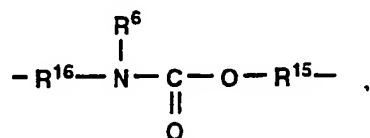
(AE)



(AF)

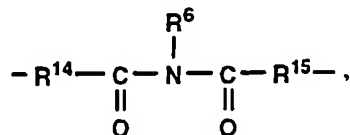


(AG)



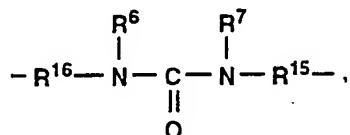
5

(AH)

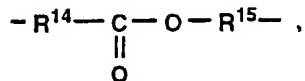


10

(AI)

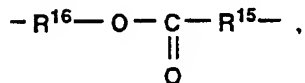


(AJ)

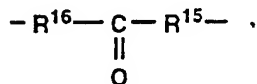


15

(AK)

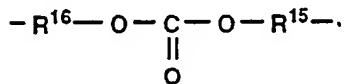


(AL)



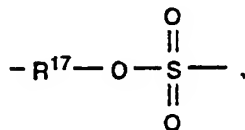
20

(AM)

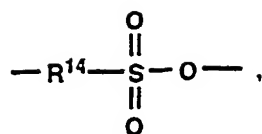


25

(AN)

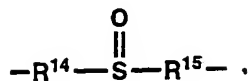


(AO)

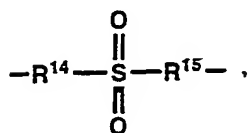


(AP)

5

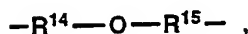


(AQ)



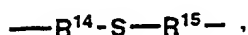
10

(AR)



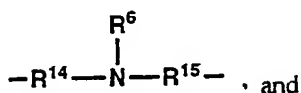
15

(AS)



(AT)

20



(AU)

25



where

30

 $\text{R}^{14}$  is selected from

a chemical bond,

C<sub>1</sub>-C<sub>8</sub>-alkyl,C<sub>3</sub>-C<sub>7</sub>-cycloalkylC<sub>2</sub>-C<sub>5</sub>-alkenyl,C<sub>3</sub>-C<sub>5</sub>-alkynyl,C<sub>6</sub>-C<sub>10</sub>-aryl,

35

C<sub>1</sub>-C<sub>3</sub>-alkyl-C<sub>6</sub>-C<sub>12</sub>-aryl,C<sub>1</sub>-C<sub>2</sub>-alkyl-C<sub>6</sub>-C<sub>10</sub>-aryl-C<sub>1</sub>-C<sub>2</sub>-alkyl,C<sub>6</sub>-C<sub>10</sub>-aryl-C<sub>1</sub>-C<sub>2</sub>-alkyl, andC<sub>6</sub>-C<sub>10</sub>-aryloxy-C<sub>1</sub>-C<sub>2</sub>-alkyl,

40

 $\text{R}^{15}$  is selected from

a chemical bond,

C<sub>1</sub>-C<sub>4</sub>-alkyl,C<sub>2</sub>-C<sub>4</sub>-alkenyl,C<sub>2</sub>-C<sub>4</sub>-alkynyl,C<sub>6</sub>-C<sub>10</sub>-aryl, and

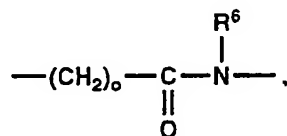
45

C<sub>1</sub>-C<sub>3</sub>-alkyl-C<sub>6</sub>-C<sub>12</sub>-aryl,

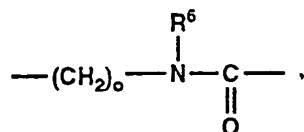
- 5  $R^{16}$  is selected from  
 a chemical bond,  
 C<sub>1</sub>-C<sub>5</sub>-alkyl,  
 C<sub>3</sub>-C<sub>7</sub>-cycloalkyl,  
 C<sub>3</sub>-C<sub>5</sub>-alkenyl,  
 C<sub>3</sub>-C<sub>5</sub>-alkynyl,  
 C<sub>6</sub>-C<sub>10</sub>-aryl,  
 C<sub>1</sub>-C<sub>3</sub>-alkyl-C<sub>6</sub>-C<sub>12</sub>-aryl, and  
 C<sub>6</sub>-C<sub>10</sub>-aryl-C<sub>1</sub>-C<sub>2</sub>-alkyl,
- 10  $R^{17}$  is selected from  
 C<sub>3</sub>-C<sub>4</sub>-alkenyl,  
 C<sub>3</sub>-C<sub>4</sub>-alkynyl,  
 C<sub>6</sub>-C<sub>10</sub>-aryl, and  
 benzyl, and
- 15 HET is a heterocycle having from 5-14 atoms in from 1-3 cycle(s), each cycle having from 1-3 heteroatoms selected from N, O, and S,  
 where the substituents are selected from one to three  $R^{12}$ ;

20  $L^2$  is an optionally substituted bivalent radical selected from the group consisting of

(A)

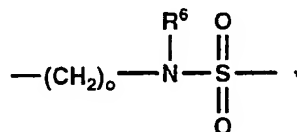


(B)



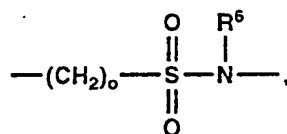
25

(C)

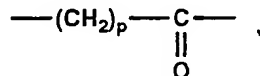


30

(D)

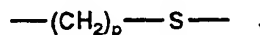


(E)

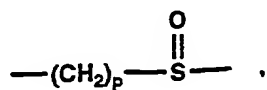


35

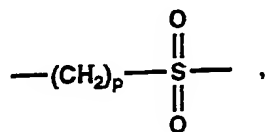
(F)



(G)

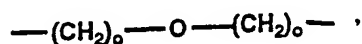


(H)



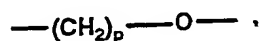
5

(I)



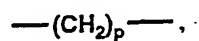
10

(J)

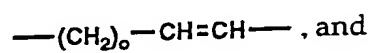


15

(K)

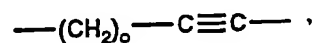


(L)



20

(M)



25

where

o is 1 or 2,

p is 1, 2, 3 or 4,

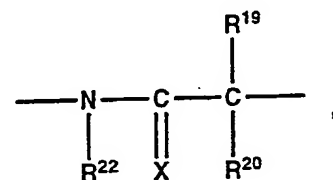
 $\text{R}^6$  is defined above, and

30

where the substituents are selected from one to three  $\text{R}^{12}$ ;

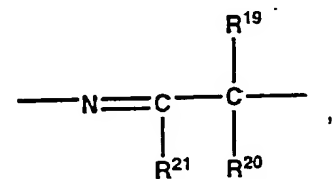
T-U-G is selected from the group consisting of

(A)



35

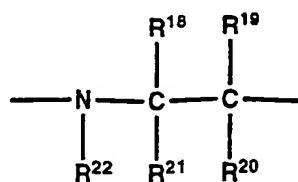
(B)



40

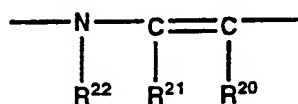


(C)

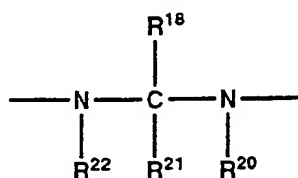


5

(D)

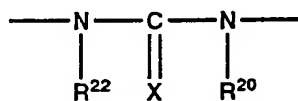


(E)



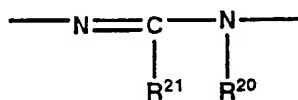
10

(F)

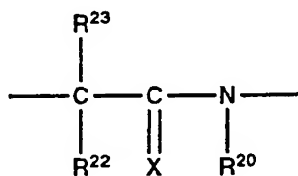


15

(G)

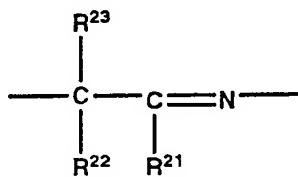


(H)



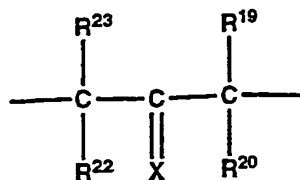
20

(I)



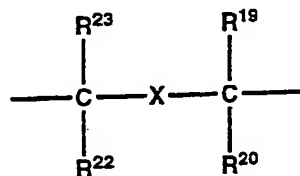
25

(J)

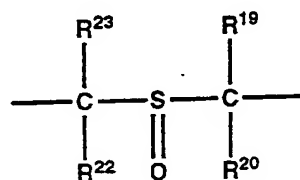


5

(K)

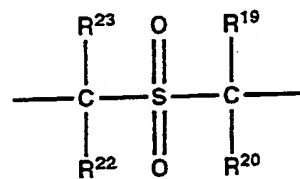


(L)



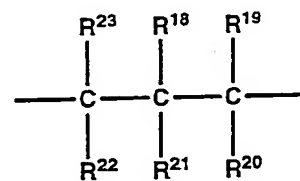
10

(M)



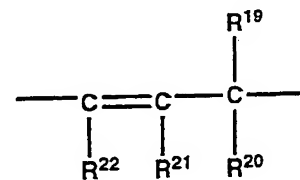
15

(N)

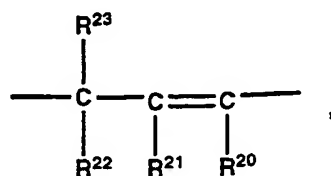


20

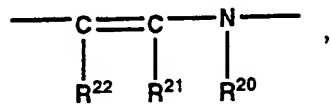
(O)



(P)

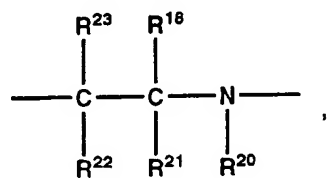


(Q)



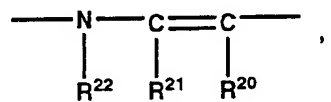
5

(R)



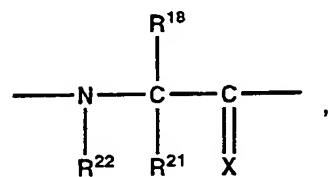
10

(S)

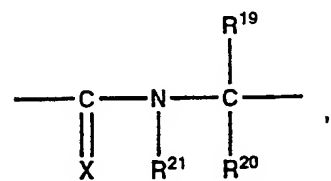


15

(T)

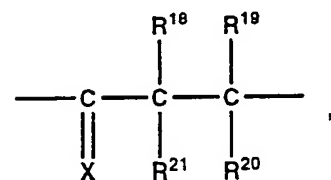


(U)



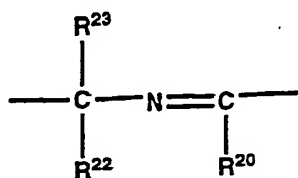
20

(V)



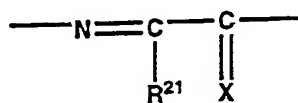
25

(W)

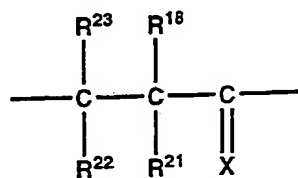


5

(X)

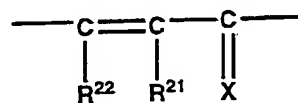


(Y)



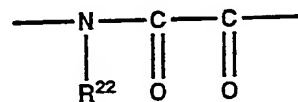
10

(Z)



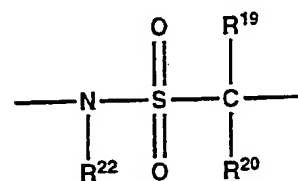
15

(AA)



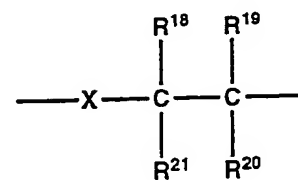
20

(AB)

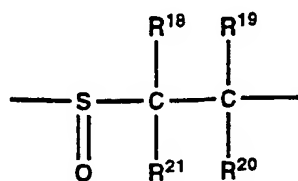


25

(AC)

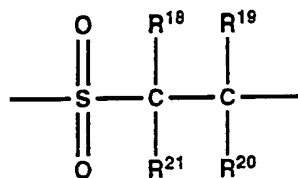


(AD)



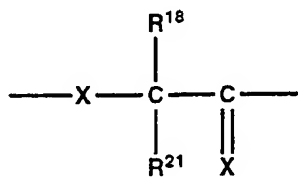
5

(AE)

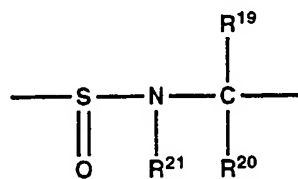


10

(AF)

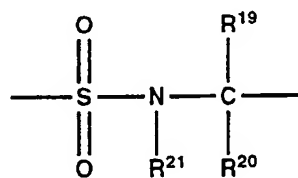


(AG)



15

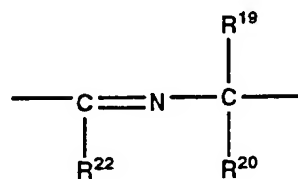
(AH)



and

20

(AI)



25

where

X is O or S;

R<sup>18</sup> and R<sup>23</sup> are independently selected from the group consisting of

- 5 (i) hydrogen,  
 (ii) optionally substituted C<sub>1</sub>-C<sub>12</sub> alkyl,  
 (iii) optionally substituted C<sub>3</sub>-C<sub>12</sub> alkenyl,  
 (iv) optionally substituted C<sub>3</sub>-C<sub>14</sub> cycloalkyl,  
 (v) optionally substituted C<sub>1</sub>-C<sub>12</sub> alkyl-C<sub>6</sub>-C<sub>14</sub>-aryl,  
 (vi) optionally substituted C<sub>6</sub>-C<sub>14</sub> aryl,  
 (vii) optionally substituted C<sub>1</sub>-C<sub>4</sub> alkylphenyl, and  
 (viii) optionally substituted C<sub>1</sub>-C<sub>12</sub> alkoxy,
- 10 R<sup>19</sup> and R<sup>20</sup> are independently selected from  
 (i) hydrogen,  
 (ii) halo(F, Cl, Br, I),  
 (iii) C<sub>1</sub>-C<sub>4</sub> alkoxy,  
 (iv) C<sub>1</sub>-C<sub>4</sub> alkyl,  
 (v) phenyl,  
 (vi) benzyl, and  
 (vii) halo(F, Cl, Br, I)-C<sub>1</sub>-C<sub>4</sub>-alkyl,
- 15 R<sup>21</sup> and R<sup>22</sup> are independently selected from the group consisting of  
 (i) hydrogen,  
 (ii) optionally substituted C<sub>1</sub>-C<sub>12</sub>-alkyl,  
 (iii) optionally substituted C<sub>6</sub>-C<sub>14</sub>-aryl,  
 (iv) optionally substituted C<sub>3</sub>-C<sub>14</sub>-cycloalkyl,  
 (v) optionally substituted C<sub>1</sub>-C<sub>12</sub>-alkyl-C<sub>6</sub>-C<sub>14</sub>-aryl, and  
 (vi) optionally substituted C<sub>1</sub>-C<sub>12</sub>-alkyl-C<sub>3</sub>-C<sub>14</sub>-cycloalkyl,  
 where the substituents are selected from
- 20 (a) halo (F, Cl, Br, I),  
 (b) nitro,  
 (c) hydroxy,  
 (d) carboxy,  
 (e) tetrazole,  
 (f) hydroxamate,  
 (g) sulfonamide,  
 (h) trifluoroimide,  
 (i) phosphonate,  
 (j) C<sub>1</sub>-C<sub>6</sub>-alkyl,  
 (k) C<sub>6</sub>-C<sub>14</sub>-aryl,  
 (l) benzyl,  
 (m) C<sub>3</sub>-C<sub>14</sub>-cycloalkyl,  
 (n) COR<sup>24</sup>
- 25 where R<sup>24</sup> is selected from the group  
 C<sub>1</sub>-C<sub>8</sub> alkoxy,  
 C<sub>3</sub>-C<sub>12</sub> alkenoxy,  
 C<sub>6</sub>-C<sub>12</sub> aryloxy,  
 C<sub>1</sub>-C<sub>6</sub>-alkyl-C<sub>6</sub>-C<sub>12</sub>-aryloxy,  
 di-C<sub>1</sub>-C<sub>8</sub>-alkylamino-C<sub>1</sub>-C<sub>8</sub>-alkoxy,  
 acylamino-C<sub>1</sub>-C<sub>8</sub>-alkoxy selected from the group  
 30 acetylaminooethoxy,  
 nicotinoylaminoethoxy, and  
 succinamidoethoxy,  
 C<sub>1</sub>-C<sub>8</sub>-alkoxyloxy-C<sub>1</sub>-C<sub>8</sub>-alkoxy, and  
 C<sub>6</sub>-C<sub>12</sub> aryl-C<sub>1</sub>-C<sub>8</sub>-alkoxy where the aryl group is  
 35 unsubstituted or substituted with one to three of  
 the groups  
 nitro,  
 halo (F, Cl, Br, I),  
 40 C<sub>1</sub>-C<sub>4</sub>-alkoxy,
- 45
- 50
- 55

amino,  
hydroxy,  
hydroxy-C<sub>2</sub>-C<sub>8</sub>-alkoxy, and  
dihydroxy-C<sub>3</sub>-C<sub>8</sub>-alkoxy, and

5

(o) CONR<sup>25</sup>R<sup>26</sup>

where R<sup>25</sup> and R<sup>26</sup> are independently selected from

hydrogen,  
C<sub>1</sub>-C<sub>10</sub>-alkyl,  
C<sub>3</sub>-C<sub>10</sub>-alkenyl,

10

C<sub>6</sub>-C<sub>14</sub>-aryl,  
C<sub>1</sub>-C<sub>6</sub>-alkyl-C<sub>6</sub>-C<sub>10</sub>-aryl,

optionally R<sup>25</sup> and R<sup>26</sup> taken together may form  
trimethylene,  
tetramethylene,  
pentamethylene, and  
3-oxopentamethylene, and

15

(vii) Q<sup>2</sup>-L<sup>3</sup>.

where Q<sup>2</sup> is selected from

hydrogen, and  
Q<sup>1</sup>, and

20

L<sup>3</sup> is selected from  
a chemical bond,  
L<sup>1</sup>, and  
L<sup>2</sup>;

25

D is selected from the group consisting of

R<sup>22</sup>, and

-(C=O)-Xaa, where Xaa is one to three D or L  $\alpha$ -amino acid residues;

30 W is -R<sup>27</sup>-w where

R<sup>27</sup> is selected from

(a) a covalent bond,  
(b) substituted or unsubstituted methylene, and  
(c) substituted or unsubstituted ethylene,

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where the substituents are selected from

(i) nitro,  
(ii) halo(F, Cl, Br, I),  
(iii) C<sub>1</sub>-C<sub>6</sub> alkyl,  
(iv) halo(F, Cl, Br, I)-C<sub>1</sub>-C<sub>6</sub> alkyl, and  
(v) substituted or unsubstituted phenyl

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where the phenyl substituents are selected from

(1) C<sub>1</sub>-C<sub>6</sub> alkyl,  
(2) C<sub>1</sub>-C<sub>6</sub> alkoxy,  
(3) halo(F, Cl, Br, I), and  
(4) CF<sub>3</sub>;

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w is selected from

(a) -COR<sup>28</sup>,  
(b) -SO<sub>3</sub>R<sup>31</sup>,  
(c) -NHSO<sub>2</sub>R<sup>32</sup>,  
(d) -PO(OR<sup>31</sup>)<sub>2</sub>,  
(e) -SO<sub>2</sub>NHR<sup>32</sup>,  
(f) -CONHOR<sup>31</sup>,  
(g) -C(OH)R<sup>33</sup>PO(OR<sup>33</sup>)<sub>2</sub>,  
(h) -CN,

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(i)  $-\text{SO}_2\text{NH}$ -heterocycle where the heterocycle is a 5- or 6-member aromatic ring containing 1 to 3 heteroatoms selected from O, N, and S and where the heterocycle is unsubstituted or substituted with one or two substituents selected from the group

- (i)  $-\text{OH}$ ,
- (ii)  $-\text{SH}$ ,
- (iii)  $-(\text{C}_1-\text{C}_4\text{alkyl})$ ,
- (iv)  $-\text{C}_1-\text{C}_4\text{alkoxyl}$ ,
- (v)  $\text{CF}_3$ ,
- (vi) halo(F, Cl, Br, I),
- (vii)  $\text{NO}_2$ ,
- (viii)  $-\text{COOH}$ ,
- (ix)  $-\text{COO}-(\text{C}_1-\text{C}_4\text{alkyl})$ ,
- (x)  $-\text{NH}_2$ ,
- (xi)  $-\text{NH}(\text{C}_1-\text{C}_4\text{alkyl})$ , and
- (xii)  $-\text{N}(\text{C}_1-\text{C}_4\text{alkyl})_2$ ,

(j)  $-\text{CH}_2\text{SO}_2\text{NH}$ -heterocycle,

(k)  $-\text{SO}_2\text{NHCOR}^{33}$ ,

(l)  $-\text{CH}_2\text{SO}_2\text{NHCOR}^{32}$ ,

(m)  $-\text{CONHSO}_2\text{R}^{33}$ ,

(n)  $-\text{CH}_2\text{CONHSO}_2\text{R}^{33}$ ,

(o)  $-\text{NHCONHSO}_2\text{R}^{33}$ ,

(p)  $-\text{NHSO}_2\text{NHCOR}^{33}$ ,

(q)  $-\text{CONHNHSO}_2\text{CF}_3$ ,

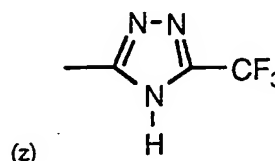
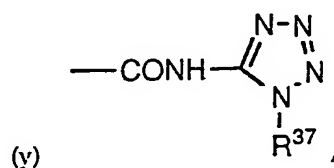
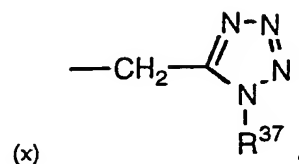
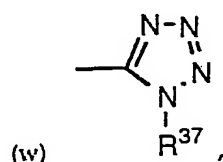
(r)  $-\text{CON}(\text{OH})\text{R}^{31}$ ,

(s)  $-\text{CONHCOCF}_3$ ,

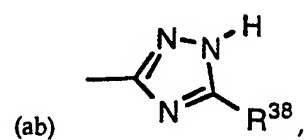
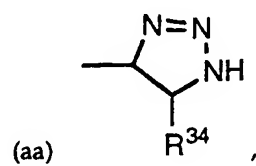
(t)  $-\text{CONHSO}_2\text{R}^{28}$ ,

(u)  $-\text{CONHSO}_2\text{R}^{29}$ ,

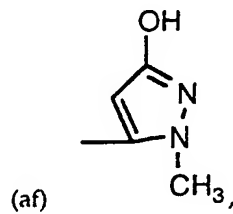
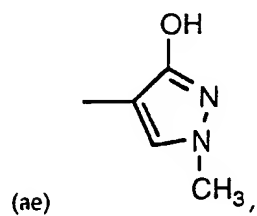
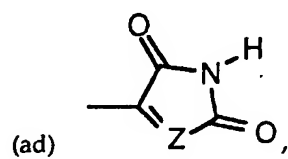
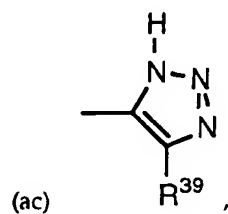
(v)  $-\text{CONHSO}_2\text{R}^{30}$ ,



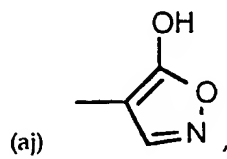
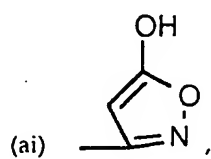
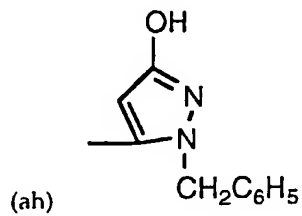
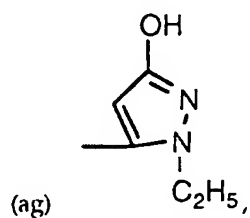




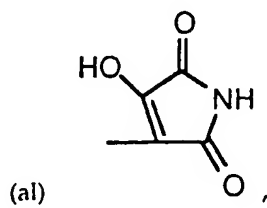
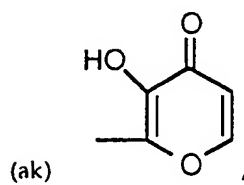
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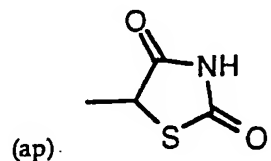
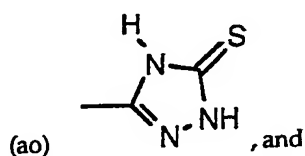
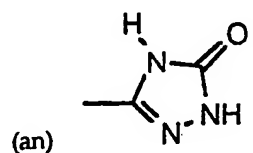
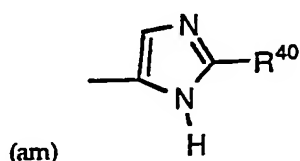


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R<sup>28</sup> is selected from the group consisting of

- (a) hydroxy,  
 (b) C<sub>1</sub>-C<sub>8</sub>-alkoxy,  
 (c) C<sub>3</sub>-C<sub>12</sub>-alkenoxy,  
 (d) C<sub>6</sub>-C<sub>12</sub>-aryloxy,  
 (e) C<sub>1</sub>-C<sub>6</sub>-alkyl-C<sub>6</sub>-C<sub>12</sub>-aryloxy,  
 (f) di-C<sub>1</sub>-C<sub>8</sub>-alkylamino-C<sub>1</sub>-C<sub>8</sub>-alkoxy,  
 (g) acylamino-C<sub>1</sub>-C<sub>8</sub>-alkoxy selected from the group  
     (i) acetylaminooethoxy,  
     (ii) nicotinoylaminoethoxy, and  
     (iii) succinamidoethoxy,  
 (h) C<sub>1</sub>-C<sub>8</sub>-alkoxyloxy-C<sub>1</sub>-C<sub>8</sub>-alkoxy,  
 (i) C<sub>6</sub>-C<sub>12</sub>-aryl-C<sub>1</sub>-C<sub>8</sub>-alkoxy,  
 where any aryl groups are unsubstituted or substituted with one to  
 three of the groups  
     (i) nitro,  
     (ii) halo (F, Cl, Br, I),  
     (iii) C<sub>1</sub>-C<sub>4</sub>-alkoxy, and  
     (iv) amino,  
 (j) hydroxy-C<sub>2</sub>-C<sub>8</sub>-alkoxy,  
 (k) dihydroxy-C<sub>3</sub>-C<sub>8</sub>-alkoxy, and  
 (l) NR<sup>29</sup>R<sup>30</sup>;

R<sup>29</sup> and R<sup>30</sup> are independently selected from the group

- (a) hydrogen,  
 (b) C<sub>1</sub>-C<sub>8</sub>-alkyl,  
 (c) C<sub>3</sub>-C<sub>8</sub>-alkenyl,  
 (d) C<sub>6</sub>-C<sub>12</sub>-aryl,  
 (e) C<sub>6</sub>-C<sub>12</sub>-aryl-C<sub>1</sub>-C<sub>8</sub>-alkyl  
 where any aryl groups are unsubstituted or substituted with one to three of  
 the groups  
     (i) nitro,  
     (ii) halo (F, Cl, Br, I),  
     (iii) C<sub>1</sub>-C<sub>4</sub>-alkoxy, and  
     (iv) amino;

R<sup>31</sup> is selected from the group consisting of

- (a) hydrogen,  
 (b) C<sub>1</sub>-C<sub>6</sub> alkyl,  
 (c) halo(F, Cl, Br, I)-C<sub>1</sub>-C<sub>6</sub> alkyl,  
 (d) phenyl,

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- (e) benzyl, and
- (f)  $-\text{CH}_2-\text{O}-\text{COCH}_3$ ;

$\text{R}^{32}$  is selected from the group consisting of

- (a) hydrogen,
- (b) benzyl and
- (c)  $-\text{CH}(\text{R}^{35})-\text{O}-\text{C}(\text{O})\text{R}^{35}$ ;

$\text{R}^{33}$  is selected from the group consisting of

- (a) aryl,
- (b) heteroaryl,
- (c)  $(\text{C}_3-\text{C}_7)$ -cycloalkyl,
- (d)  $(\text{C}_1-\text{C}_4)$ -alkyl, unsubstituted or substituted with a substituent selected from the group consisting of

- (i) aryl,
- (ii) heteroaryl,
- (iii)  $-\text{OH}$ ,
- (iv)  $-\text{SH}$ ,
- (v)  $(\text{C}_1-\text{C}_4)$ -alkyl,
- (vi)  $(\text{C}_1-\text{C}_4)$ -alkoxy,
- (vii)  $(\text{C}_1-\text{C}_4)$ -alkylthio,
- (viii)  $-\text{CF}_3$ ,
- (ix) halo (F, Cl, Br, I),
- (x)  $-\text{NO}_2$ ,
- (xi)  $-\text{CO}_2\text{H}$ ,
- (xii)  $\text{CO}_2-(\text{C}_1-\text{C}_4)$ -alkyl,
- (xiii)  $-\text{NH}_2$ ,
- (xiv)  $-\text{N}[(\text{C}_1-\text{C}_4)\text{-alkyl}]_2$ ,
- (xv)  $-\text{NH}[(\text{C}_1-\text{C}_4)\text{-alkyl}]$ ,
- (xvi)  $-\text{PO}_3\text{H}$ , and
- (xvii)  $\text{PO}(\text{OH})(\text{C}_1-\text{C}_4)\text{-alkoxy}$ , and

- (e)  $(\text{C}_1-\text{C}_4)$ -perfluoroalkyl;

$\text{R}^{34}$  is selected from the group consisting of

- (a)  $-\text{CN}$ ,
- (b)  $-\text{NO}_2$ ,
- (c)  $-\text{COOR}^{31}$ ,
- (d)  $\text{C}_1-\text{C}_6$ -perfluoroalkyl, and
- (e)  $\text{CF}_3$ ;

$\text{R}^{35}$  is independently selected from the group consisting of

- (a) hydrogen,
  - (b) optionally substituted  $(\text{C}_1-\text{C}_6)$ -alkyl,
  - (c) optionally substituted  $(\text{C}_2-\text{C}_6)$ -alkenyl,
  - (d) optionally substituted  $(\text{C}_2-\text{C}_6)$ -alkynyl, and
  - (e) optionally substituted  $(\text{C}_3-\text{C}_8)$ -cycloalkyl,
- where the substituents are selected from the group

- (i)  $\text{OH}$ ,
- (ii)  $(\text{C}_1-\text{C}_4)$ -alkoxy,
- (iii)  $\text{CO}_2\text{R}^{33}$ ,
- (iv)  $\text{OCOR}^{33}$ ,
- (v)  $\text{CONHR}^{33}$ ,
- (vi)  $\text{CON}(\text{R}^{33})_2$ ,
- (vii)  $\text{N}(\text{R}^{33})\text{C}(\text{O})\text{R}^{33}$ ,
- (viii)  $\text{NH}_2$ ,
- (ix)  $(\text{C}_1-\text{C}_4)$ -alkylamino,
- (x)  $\text{di}[(\text{C}_1-\text{C}_4)\text{-alkyl}]\text{amino}$ ,
- (xi) aryl, and

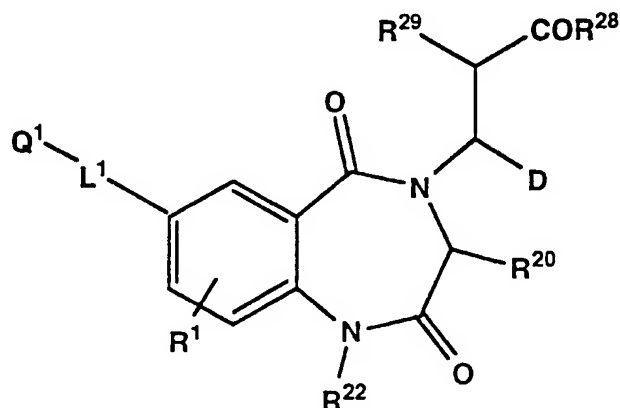
- (xii) heteroaryl,
- (f) -C(O)-aryl,
- (g) -NO<sub>2</sub>,
- (h) halo(Cl, Br, I, F),
- (i) -OH,
- (j) -OR<sup>36</sup>,
- (k) -(C<sub>1</sub>-C<sub>4</sub>)-perfluoroalkyl,
- (l) -SH,
- (m) -S(O)<sub>1-2</sub> (C<sub>1</sub>-C<sub>4</sub>)-alkyl,
- (n) -CO<sub>2</sub>R<sup>33</sup>,
- (o) -SO<sub>3</sub>H,
- (p) -NR<sup>33</sup>R<sup>36</sup>,
- (q) -NR<sup>33</sup>C(O)R<sup>36</sup>,
- (r) -NR<sup>33</sup>COOR<sup>32</sup>,
- (s) -SO<sub>2</sub>NHR<sup>32</sup>,
- (t) -SO<sub>2</sub>NR<sup>33</sup>R<sup>33</sup>,
- (u) -NHSO<sub>2</sub>R<sup>32</sup>,
- (v) -C(O)NHSO<sub>2</sub>R<sup>32</sup>,
- (w) aryl,
- (x) heterocycle,
- (y) morpholin-4-yl,
- (z) CONH<sub>2</sub>, and
- (aa) 1H-tetrazol-5-yl;
- R<sup>36</sup> is selected from the group
- (a) hydrogen, and
- (b) (C<sub>1</sub>-C<sub>4</sub>)-alkyl unsubstituted or substituted with
- (i) NH<sub>2</sub>,
- (ii) NH[(C<sub>1</sub>-C<sub>4</sub>)-alkyl],
- (iii) N[(C<sub>1</sub>-C<sub>4</sub>)-alkyl]<sub>2</sub>,
- (iv) CO<sub>2</sub>H,
- (v) CO<sub>2</sub>(C<sub>1</sub>-C<sub>4</sub>)-alkyl,
- (vi) OH,
- (vii) SO<sub>3</sub>H, and
- (viii) SO<sub>2</sub>NH<sub>2</sub>;
- R<sup>37</sup> is selected from the group consisting of
- (a) hydrogen,
- (b) (C<sub>1</sub>-C<sub>6</sub>)-alkyl,
- (c) (C<sub>2</sub>-C<sub>6</sub>)-alkenyl,
- (d) (C<sub>1</sub>-C<sub>6</sub>)-alkoxyalkyl,
- (e) -CH<sub>2</sub>-O-COCH<sub>3</sub>, and
- (f) -CH<sub>2</sub>-phenyl, where the phenyl is unsubstituted or substituted with a substituent selected from
- (i) -NO<sub>2</sub>,
- (ii) -NH<sub>2</sub>,
- (iii) -OH, and
- (iv) -OCH<sub>3</sub>;
- R<sup>38</sup>, R<sup>39</sup>, and R<sup>40</sup> are each independently selected from
- (a) hydrogen,
- (b) Cl,
- (c) CN,
- (d) NO<sub>2</sub>,

- (e) CF<sub>3</sub>,
- (f) C<sub>2</sub>F<sub>5</sub>,
- (g) C<sub>3</sub>F<sub>7</sub>,
- (h) CHF<sub>2</sub>,
- (i) CH<sub>2</sub>F,
- (j) CO<sub>2</sub>CH<sub>3</sub>,
- (k) CO<sub>2</sub>C<sub>2</sub>H<sub>5</sub>,
- (l) SO<sub>2</sub>CH<sub>3</sub>,
- (m) SO<sub>2</sub>CF<sub>3</sub>, and
- (n) SO<sub>2</sub>C<sub>6</sub>F<sub>5</sub>,

wherein Z is selected from O, S, NR<sup>41</sup> and CH<sub>2</sub>;  
 R<sup>41</sup> is selected from hydrogen, CH<sub>3</sub>, and CH<sub>2</sub>C<sub>6</sub>H<sub>5</sub>; and

pharmaceutically acceptable salts thereof.

2. The compound of Claim 1 represented by formula II:



(II)

where

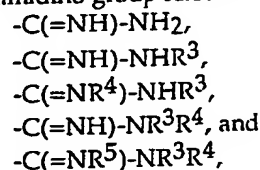
- 25 R<sup>1</sup> is one to three groups independently selected from hydrogen, halo(F, Cl, Br, I), cyano, carboxamido, carbamoyloxy, formyloxy, formyl, azido, nitro, ureido, thioureido, hydroxy, mercapto, sulfonamido, and an optionally substituted radical selected from C<sub>1</sub>-C<sub>12</sub> alkyl, C<sub>2</sub>-C<sub>12</sub> alkenyl, C<sub>3</sub>-C<sub>12</sub> alkynyl, C<sub>3</sub>-C<sub>12</sub> cycloalkyl, C<sub>6</sub>-C<sub>14</sub> aryl, C<sub>6</sub>-C<sub>10</sub> aryl-C<sub>1</sub>-C<sub>8</sub> alkyl, C<sub>1</sub>-C<sub>12</sub> alkyloxy, C<sub>6</sub>-C<sub>14</sub> aryloxy, and C<sub>1</sub>-C<sub>12</sub> acylamino, where the substituents are selected from halo(F, Cl, Br, I), cyano, azido, nitro, hydroxy, mercapto, sulfonamido, ureido, thioureido, carboxamido, carbamoyloxy, formyloxy, formyl, C<sub>1</sub>-C<sub>4</sub> alkyl, C<sub>1</sub>-C<sub>4</sub> alkoxy, phenyl, and phenoxy;

- 35 Q<sup>1</sup> is selected from the group consisting of  
 (A) an amino group selected from  
 -NH<sub>2</sub>,  
 -NR<sup>3H</sup>,  
 -NR<sup>3R4</sup>, and

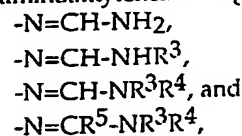


where  $\text{R}^3$ ,  $\text{R}^4$ , and  $\text{R}^5$  are independently selected from (i) an optionally substituted radical selected from (a)  $-\text{NR}^6\text{R}^7$ , (b)  $-\text{C}(=\text{NR}^8)-\text{NR}^6\text{R}^7$ , (c)  $\text{N}=\text{CR}^9-\text{NR}^6\text{R}^7$ , (d)  $-\text{NR}^{10}-\text{CR}^9=\text{NR}^8$ , and (e)  $-\text{NR}^{10}-\text{C}(=\text{NR}^8)-\text{NR}^6\text{R}^7$

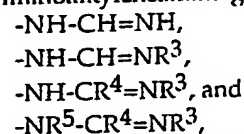
- 5 where each  $\text{R}^6$ ,  $\text{R}^7$ ,  $\text{R}^8$ ,  $\text{R}^9$ , and  $\text{R}^{10}$  is independently selected from hydrogen, C<sub>1</sub>-C<sub>4</sub> alkoxy, C<sub>1</sub>-C<sub>4</sub> alkyl, and halo(F, Cl, Br, I)-C<sub>1</sub>-C<sub>4</sub>-alkyl, (ii) optionally substituted C<sub>1</sub>-C<sub>12</sub> alkyl, (iii) optionally substituted C<sub>3</sub>-C<sub>12</sub> cycloalkyl, (iv) optionally substituted C<sub>6</sub>-C<sub>14</sub> aryl, (v) optionally substituted C<sub>1</sub>-C<sub>6</sub> alkyl-C<sub>6</sub>-C<sub>14</sub>-aryl, (vi) optionally substituted C<sub>1</sub>-C<sub>8</sub> alkoxy, and (vii) optionally substituted C<sub>6</sub>-C<sub>14</sub> aryloxy, where the substituents are one to three  $\text{R}^{11}$ , each  $\text{R}^{11}$  independently selected from (a) optionally substituted C<sub>6</sub>-C<sub>12</sub> aryloxy, (b) optionally substituted C<sub>6</sub>-C<sub>12</sub> arylamino, (c) optionally substituted C<sub>6</sub>-C<sub>12</sub> aroyl, where the substituents are one to three  $\text{R}^{12}$ , each  $\text{R}^{12}$  independently selected from nitro, amino, C<sub>1</sub>-C<sub>8</sub> alkylamino, di-(C<sub>1</sub>-C<sub>8</sub>) alkylamino, amidino, aminomethyleneimino, imino, imino-C<sub>1</sub>-C<sub>4</sub> alkyl, iminomethyleneamino, guanidino, C<sub>6</sub>-C<sub>10</sub> arylamino, C<sub>1</sub>-C<sub>8</sub> acylamino, C<sub>1</sub>-C<sub>4</sub> alkylsulfonamino, azido, cyano, hydroxy, hydroxy-C<sub>1</sub>-C<sub>8</sub>-alkyl, C<sub>1</sub>-C<sub>8</sub>-alkoxy, phenyloxy, C<sub>1</sub>-C<sub>8</sub> alkanoyloxy, C<sub>1</sub>-C<sub>8</sub> alkanoyl, C<sub>6</sub>-C<sub>12</sub> aroyl, benzamido, phenyl, halo(F, Cl, Br, I), halo-C<sub>1</sub>-C<sub>8</sub>-alkyl, and C<sub>1</sub>-C<sub>8</sub>-alkyl, (d) halo(F, Cl, Br, I), (e) hydroxy, (f) mercapto, (g) formyl, (h) formyloxy, (i) carboxy, (j) amino, (k) ureido, (l) amidino, (m) guanidino, (n) aminomethylimino, (o) imino, (p) glycol, (q) phthalimido, (r) succinimido, (s) morpholino, and (t) C<sub>3</sub>-C<sub>7</sub> cycloalkyl, optionally  $\text{R}^3$  and  $\text{R}^4$  taken together may form optionally substituted tetramethylene, pentamethylene, 3-oxopentamethylene, and 3-azapentamethylene, where the substituents are selected from one to three  $\text{R}^{12}$ ,  
 25 (B) an amidino group selected from



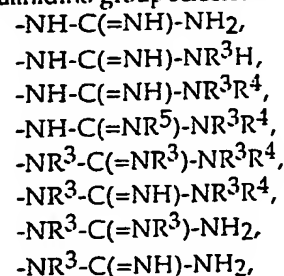
- 30 where  $\text{R}^3$ ,  $\text{R}^4$ , and  $\text{R}^5$  are defined above,  
 (C) an aminoalkyleneamino group selected from



- 35 where  $\text{R}^3$ ,  $\text{R}^4$ , and  $\text{R}^5$  are defined above,  
 (D) an iminoalkyleneamino group selected from



- 40 where  $\text{R}^3$ ,  $\text{R}^4$ , and  $\text{R}^5$  are defined above,  
 (E) a guanidino group selected from



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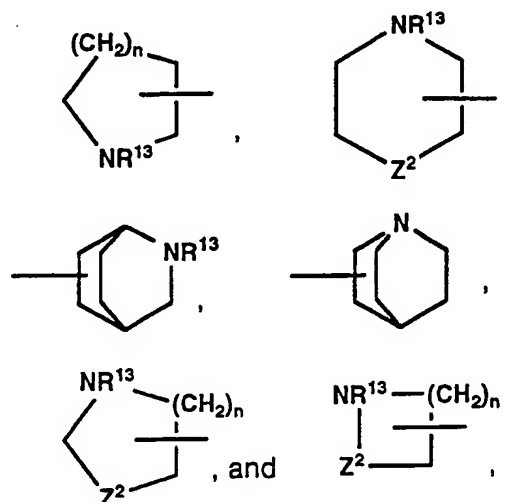
$-\text{NR}^3-\text{C}(=\text{NR}^3)-\text{NHR}^4$ , and

$-\text{NR}^3-\text{C}(=\text{NH})-\text{NHR}^4$ ,

where  $\text{R}^3$ ,  $\text{R}^4$ , and  $\text{R}^5$  are defined above, and

(F) an optionally substituted saturated heterocyclic group selected from

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where (1)  $n$  is 0, 1, 2, or 3, (2)  $\text{R}^{13}$  is selected from;  $\text{R}^6$ ,  $-\text{CR}^9=\text{NR}^8$ ,  $-\text{CR}^9(=\text{NR}^8)-\text{NR}^6\text{R}^7$ ,  $-\text{C}(=\text{NR}^8)-\text{NR}^6\text{R}^7$ ,  $-\text{N}=\text{CR}^9-\text{NR}^6\text{R}^7$ ,  $-\text{NR}^{10}-\text{CR}^9=\text{NR}^8$ , and  $-\text{NR}^{10}$ .

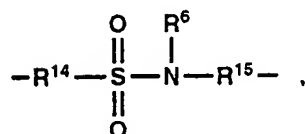
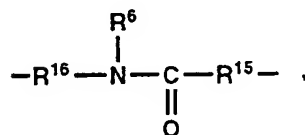
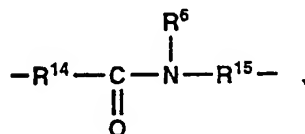
$(\text{C}=\text{NR}^8)-\text{NR}^6\text{R}^7$  where  $\text{R}^6$ ,  $\text{R}^{10}$  are defined above, (3)  $\text{Z}^2$  is O, S, or  $\text{NR}^{13}$ , and (4) the substituents are independently one to three  $\text{R}^{12}$ ;

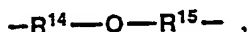
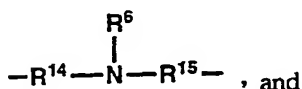
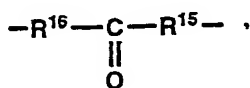
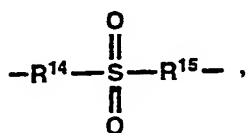
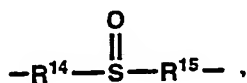
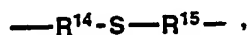
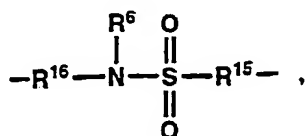
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$\text{L}^1$  is an optionally substituted bivalent radical selected from the group consisting of C3-C7-alkylene, C3-C7-cycloalkylene, C3-C7-alkenylene, C4-C7-cycloalkenylene, C5-C8-cycloalkadienylene, C3-C7-alkadienylene, C3-C7-alkynylene, C4-C7-alkenynylene, C6-C14-arylene, C6-C14-aryl-C2-C4-alkynylene, C1-C3-alkyl-C6-C14-aryl-C2-C4-alkynylene, C6-C14-aryl-C2-C4-alkenylene, C1-C3-alkyl-C6-C14-arylene, C1-C3-alkyl-C6-C14-aryl-C2-C4-alkenylene, C6-C14-aryl-C1-C3-alkylene, C6-C14-aryl-C1-C3-alkoxy, C1-C2-alkyl-C6-C14-aryl-C1-C2-alkylene, C1-C3-alkoxy-C6-C14-arylene, C2-C8-alkoxy, C1-C5-alkoxy-C1-C5-alkylene, C6-C10-aryloxy, C6-C10-aryloxy-C1-C5-alkylene, C6-C10-arylthio-C1-C5-alkylene,

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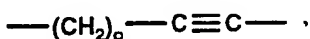
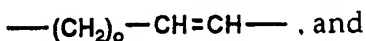
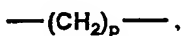
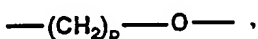
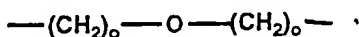
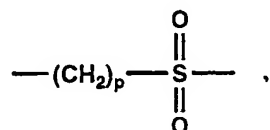
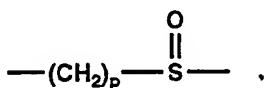
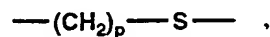
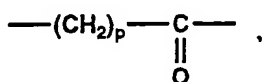
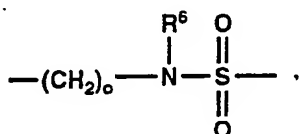
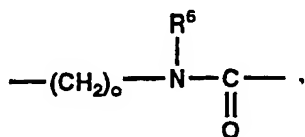
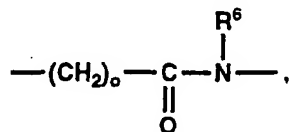
where  $\text{R}^{14}$  is selected from a chemical bond, C<sub>1</sub>-C<sub>8</sub>-alkyl, C<sub>3</sub>-C<sub>7</sub>-cycloalkyl, C<sub>2</sub>-C<sub>5</sub>-alkenyl, C<sub>3</sub>-C<sub>5</sub>-alkynyl, C<sub>6</sub>-C<sub>10</sub>-aryl, C<sub>1</sub>-C<sub>3</sub>-alkyl-C<sub>6</sub>-C<sub>12</sub>-aryl, C<sub>1</sub>-C<sub>2</sub>-alkyl-C<sub>6</sub>-C<sub>10</sub>-aryl-C<sub>1</sub>-C<sub>2</sub>-alkyl, and C<sub>6</sub>-C<sub>10</sub>-aryl-C<sub>1</sub>-C<sub>2</sub>-alkyl,  $\text{R}^{15}$  is selected from a chemical bond, C<sub>1</sub>-C<sub>4</sub>-alkyl, C<sub>2</sub>-C<sub>4</sub>-alkenyl, C<sub>2</sub>-C<sub>4</sub>-alkynyl, and C<sub>6</sub>-C<sub>10</sub>-aryl,  $\text{R}^{16}$  is selected from a chemical bond, C<sub>1</sub>-C<sub>5</sub>-alkyl, C<sub>3</sub>-C<sub>7</sub>-cycloalkyl, C<sub>3</sub>-C<sub>5</sub>-alkenyl, C<sub>3</sub>-C<sub>5</sub>-alkynyl, C<sub>6</sub>-C<sub>10</sub>-aryl, C<sub>1</sub>-C<sub>3</sub>-alkyl-C<sub>6</sub>-C<sub>12</sub>-aryl, C<sub>1</sub>-C<sub>3</sub>-alkyl-C<sub>6</sub>-C<sub>12</sub>-aryl, and C<sub>6</sub>-C<sub>10</sub>-aryl-C<sub>1</sub>-C<sub>2</sub>-alkyl, where the substituents are selected from one to three  $\text{R}^{12}$ ;

$\text{R}^{20}$  is selected from the group hydrogen, halo(F, Cl, Br, I), C<sub>1</sub>-C<sub>4</sub> alkoxy, C<sub>1</sub>-C<sub>4</sub> -alkyl, phenyl, benzyl, and halo(F, Cl, Br, I)-C<sub>1</sub>-C<sub>4</sub>-alkyl;

$\text{R}^{22}$  is selected from the group consisting of (i) hydrogen, (ii) optionally substituted C<sub>1</sub>-C<sub>12</sub>-alkyl, (iii) optionally substituted C<sub>6</sub>-C<sub>14</sub>-aryl, (iv) optionally substituted C<sub>3</sub>-C<sub>14</sub>-cycloalkyl, (v) optionally substituted C<sub>1</sub>-C<sub>12</sub>-alkyl-C<sub>6</sub>-C<sub>14</sub>-aryl, (vi) optionally substituted C<sub>1</sub>-C<sub>12</sub>-alkyl-C<sub>3</sub>-C<sub>14</sub>-cycloalkyl, where the substituents are selected from (a) halo (F, Cl, Br, I), (b) nitro, (c) hydroxy, (d) carboxy, (e) tetrazole, (f) hydroxamate, (g) sulfonamide, (h) trifluoroimide, (i) phosphonate, (j) C<sub>1</sub>-C<sub>6</sub>-alkyl, (k) C<sub>6</sub>-C<sub>14</sub>-aryl, (l) benzyl, (m) C<sub>3</sub>-C<sub>14</sub>-cycloalkyl, (n) COR<sup>24</sup> where  $\text{R}^{24}$  is selected from the group C<sub>1</sub>-C<sub>8</sub>-alkoxy, C<sub>3</sub>-C<sub>12</sub>-alkenoxy, C<sub>6</sub>-C<sub>12</sub>-aryloxy, di-C<sub>1</sub>-C<sub>8</sub>-alkylamino-C<sub>1</sub>-C<sub>8</sub>-alkoxy, acylamino-C<sub>1</sub>-C<sub>8</sub>-alkoxy selected from the group acetylaminoethoxy, nicotinoylaminoethoxy, succinamidoethoxy, and pivaloyloxyethoxy, and C<sub>6</sub>-C<sub>12</sub>-aryl-C<sub>1</sub>-C<sub>8</sub>-alkoxy where the aryl group is unsubstituted or substituted with one to three of the groups nitro, halo (F, Cl, Br, I), C<sub>1</sub>-C<sub>4</sub>-alkoxy, amino, hydroxy, hydroxy-C<sub>2</sub>-C<sub>8</sub>-alkoxy, and dihydroxy-C<sub>3</sub>-C<sub>8</sub>-alkoxy, (o) CONR<sup>25</sup>R<sup>26</sup> where  $\text{R}^{25}$  and  $\text{R}^{26}$  are independently selected from hydrogen, C<sub>1</sub>-C<sub>10</sub>-alkyl, C<sub>3</sub>-C<sub>10</sub>-alkenyl, C<sub>6</sub>-C<sub>14</sub>-aryl, C<sub>1</sub>-C<sub>6</sub>-alkyl-C<sub>6</sub>-C<sub>10</sub>-aryl,



optionally R<sup>25</sup> and R<sup>26</sup> taken together may form trimethylene, tetramethylene, pentamethylene, and 3-oxopentamethylene, and (vii) Q<sup>2</sup>-L<sup>3</sup>, where Q<sup>2</sup> is selected from hydrogen, and Q<sup>1</sup>, and L<sup>3</sup> is selected from a chemical bond, L<sup>1</sup>, and L<sup>2</sup>, where L<sup>2</sup> is an optionally substituted bivalent radical selected from the group consisting of



where o is 1 or 2, p is 1, 2, 3 or 4, R<sup>6</sup> is defined above, and where the substituents are selected from one to three R<sup>12</sup>;

D is selected from the group consisting of R<sup>22</sup>, and -(C=O)-Xaa, where Xaa is one to three D or L  $\alpha$ -amino acid;

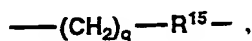
5 R<sup>28</sup> is selected from the group consisting of (i) hydroxy, (ii) C<sub>1</sub>-C<sub>8</sub>-alkoxy, (iii) C<sub>3</sub>-C<sub>12</sub>-alkenoxy, (iv) C<sub>6</sub>-C<sub>12</sub>-aryloxy, (v) C<sub>1</sub>-C<sub>6</sub>-alkyl-C<sub>6</sub>-C<sub>12</sub>-aryloxy, (vi) di-C<sub>1</sub>-C<sub>8</sub>-alkylamino-C<sub>1</sub>-C<sub>8</sub>-alkoxy, (vii) acylamino-C<sub>1</sub>-C<sub>8</sub>-alkoxy selected from the group  
 10 (a) acetylaminoethoxy, (b) nicotinoylaminoethoxy, and (c) succinamidoethoxy, (viii) C<sub>1</sub>-C<sub>8</sub>-alkoxy-C<sub>1</sub>-C<sub>8</sub>-alkoxy, (ix) C<sub>6</sub>-C<sub>12</sub>-aryl-C<sub>1</sub>-C<sub>8</sub>-alkoxy where the aryl group is unsubstituted or substituted with one to three of the groups (a) nitro, (b) halo (F, Cl, Br, I), (c) C<sub>1</sub>-C<sub>4</sub>-alkoxy, and (d) amino, (x) hydroxy-C<sub>2</sub>-C<sub>8</sub>-alkoxy, (xi) dihydroxy-C<sub>3</sub>-C<sub>8</sub>-alkoxy, and (xii) NR<sup>29</sup>R<sup>30</sup> where R<sup>29</sup> and R<sup>30</sup> are independently  
 15 selected from the group (a) hydrogen, (b) C<sub>1</sub>-C<sub>8</sub>-alkyl, (c) C<sub>3</sub>-C<sub>8</sub>-alkenyl, (d) C<sub>6</sub>-C<sub>12</sub>-aryl where the aryl group is unsubstituted or substituted with one to three of the groups nitro, halo (F, Cl, Br, I), C<sub>1</sub>-C<sub>4</sub>-alkoxy, and amino, and (e) C<sub>6</sub>-C<sub>12</sub>-aryl-C<sub>1</sub>-C<sub>8</sub>-alkyl where the aryl group is unsubstituted or substituted with one to three of the groups nitro, halo (F, Cl, Br, I), and C<sub>1</sub>-C<sub>4</sub>-alkoxy;

20 R<sup>29</sup> is one or two groups independently selected from the group (a) hydrogen, (b) C<sub>1</sub>-C<sub>8</sub>-alkyl, (c) C<sub>3</sub>-C<sub>8</sub>-alkenyl, (d) C<sub>6</sub>-C<sub>12</sub>-aryl, and (e) C<sub>6</sub>-C<sub>12</sub>-aryl-C<sub>1</sub>-C<sub>8</sub>-alkyl, where any aryl groups are unsubstituted or substituted with one to three of the groups (i) nitro, (ii) halo (F, Cl, Br, I), (iii) C<sub>1</sub>-C<sub>4</sub>-alkoxy, and (iv) amino; and

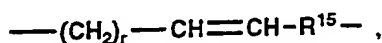
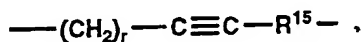
pharmaceutically acceptable salts thereof.

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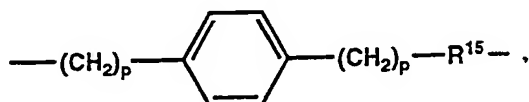
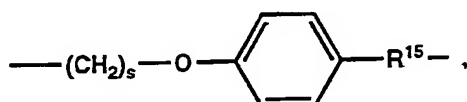
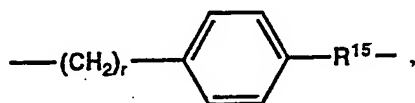
3. The compound of Claim 2 where -L<sup>1</sup>- is selected from the group consisting of



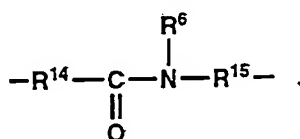
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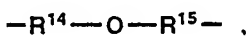
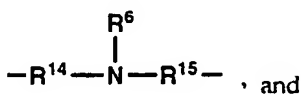
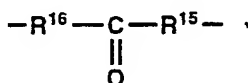
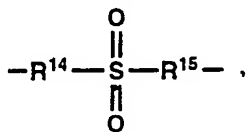
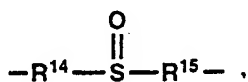
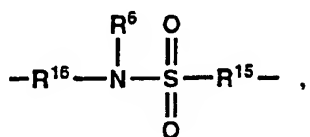
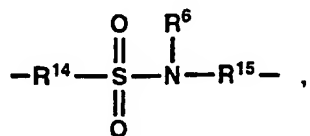
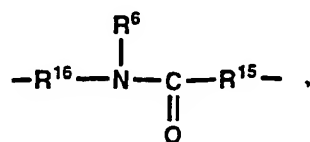


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where

p is 1, 2, 3, or 4;

q is 3, 4, 5, 6 or 7;

r is 1, 2, 3, 4, or 5;

s is 2 or 3;

R<sup>14</sup> is selected from

C<sub>2</sub>-C<sub>5</sub>-alkyl,

C<sub>3</sub>-C<sub>7</sub>-cycloalkyl,

C<sub>2</sub>-C<sub>5</sub>-alkenyl,

C<sub>3</sub>-C<sub>5</sub>-alkynyl,

C<sub>6</sub>-C<sub>10</sub>-aryl,

C<sub>1</sub>-C<sub>2</sub>-alkyl-C<sub>6</sub>-C<sub>12</sub>-aryl,

C<sub>1</sub>-C<sub>2</sub>-alkyl-C<sub>6</sub>-C<sub>10</sub>-aryl-C<sub>1</sub>-C<sub>2</sub>-alkyl,

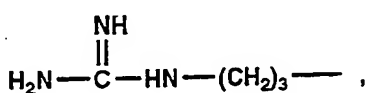
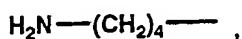
C<sub>6</sub>-C<sub>10</sub>-aryl-C<sub>1</sub>-C<sub>2</sub>-alkyl, and  
C<sub>6</sub>-C<sub>10</sub>-aryloxy-C<sub>1</sub>-C<sub>2</sub>-alkyl;

R<sup>15</sup> is a chemical bond connecting L<sup>I</sup> to position 7 of the benzodiazepinedione;

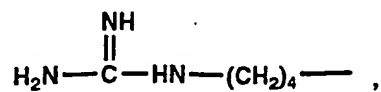
R<sup>16</sup> is selected from  
C<sub>2</sub>-C<sub>5</sub>-alkyl,  
C<sub>3</sub>-C<sub>7</sub>-cycloalkyl,  
C<sub>3</sub>-C<sub>5</sub>-alkenyl,  
C<sub>3</sub>-C<sub>5</sub>-alkynyl,  
C<sub>6</sub>-C<sub>10</sub>-aryl,  
C<sub>1</sub>-C<sub>3</sub>-alkyl-C<sub>6</sub>-C<sub>12</sub>-aryl, and  
C<sub>6</sub>-C<sub>10</sub>-aryl-C<sub>1</sub>-C<sub>2</sub>-alkyl,  
wherein R<sup>14</sup> and R<sup>16</sup> bond L<sup>1</sup> to Q<sup>1</sup>.

4. The compound of Claim 3 where R<sup>22</sup> is selected from the group consisting of  
hydrogen,  
optionally substituted C<sub>1</sub>-C<sub>6</sub>-alkyl,  
where the substituents are selected from  
amino,  
hydroxy,  
halo(F, Cl, Br, I),  
carboxy, and  
C<sub>1</sub>-C<sub>4</sub>-alkoxycarbonyl,  
optionally substituted C<sub>6</sub>-C<sub>10</sub>-aryl, and  
optionally substituted C<sub>6</sub>-C<sub>12</sub>-aryl-C<sub>1</sub>-C<sub>4</sub>-alkyl,  
where the substituents on any aryl group are selected from  
amino,  
nitro,  
halo(F, Cl, Br, I),  
halo(F, Cl, Br, I)-C<sub>1</sub>-C<sub>6</sub>-alkyl,  
C<sub>1</sub>-C<sub>6</sub>-alkoxy,  
C<sub>6</sub>-C<sub>10</sub>-aryloxy,  
amino-C<sub>1</sub>-C<sub>6</sub>-acylamino,  
amino-C<sub>1</sub>-C<sub>6</sub>-acyl, and  
guanidino-C<sub>6</sub>-C<sub>10</sub>-aroylamino.

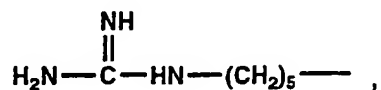
5. The compound of Claim 3 where Q<sup>1</sup>-L<sup>1</sup>- is selected from the group consisting of



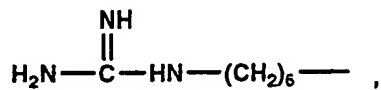
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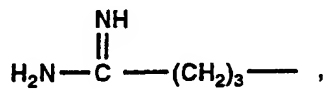
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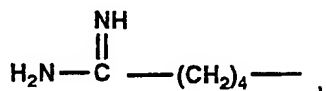
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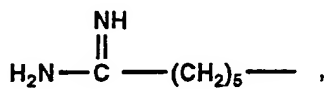
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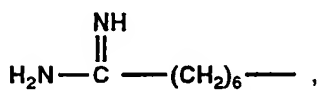
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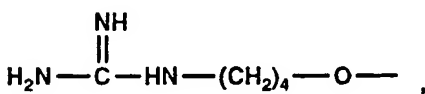
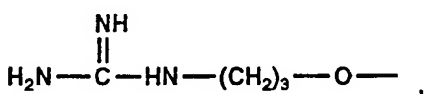
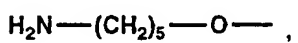
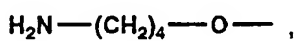
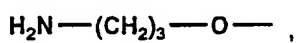
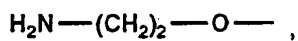
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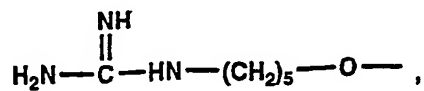


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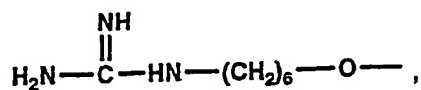


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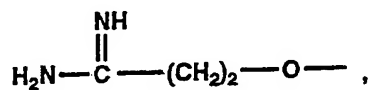




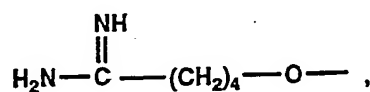
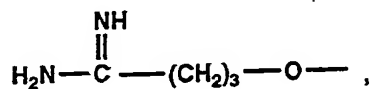
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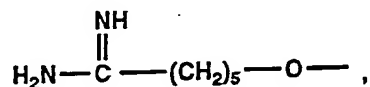
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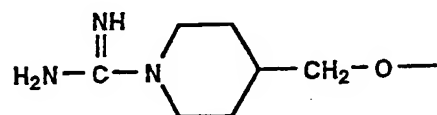
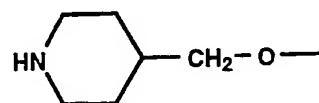
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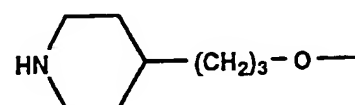
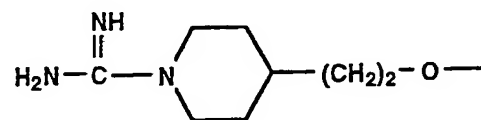
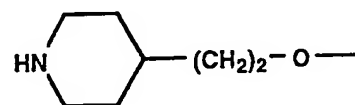
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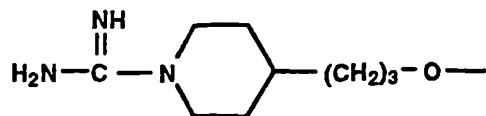


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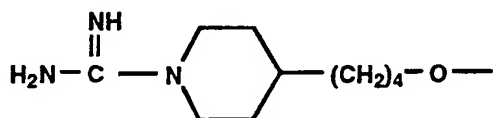
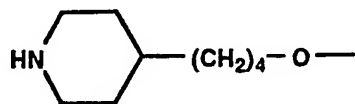


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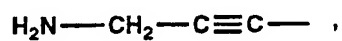




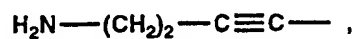
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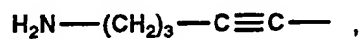
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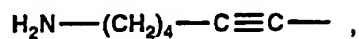
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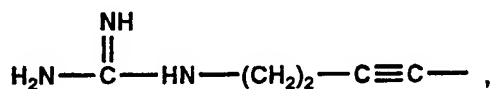
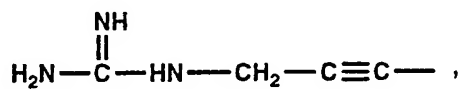
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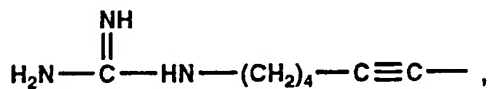
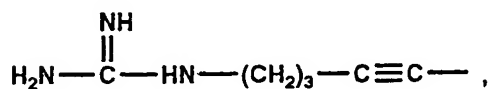
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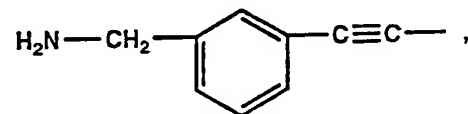
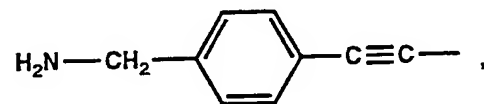
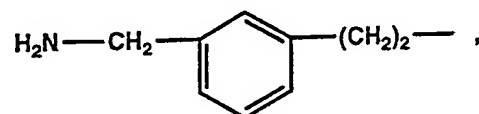
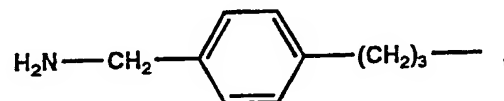
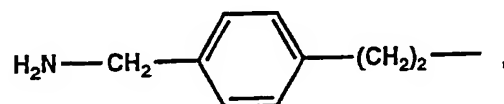
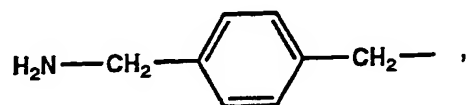
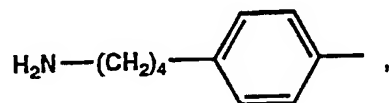
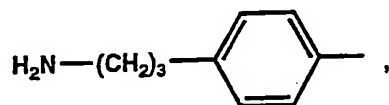
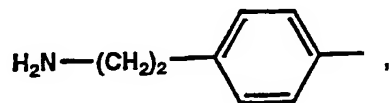
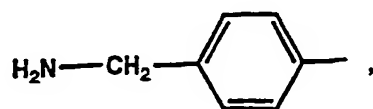


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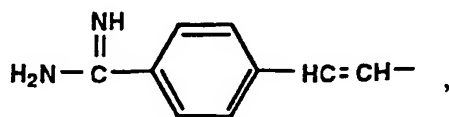
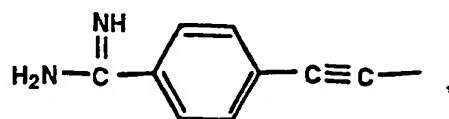


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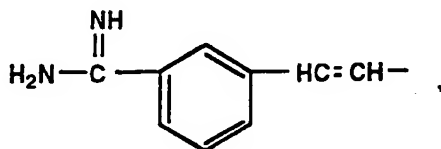




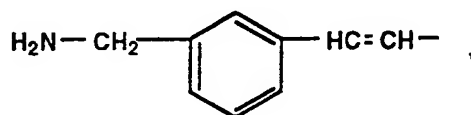




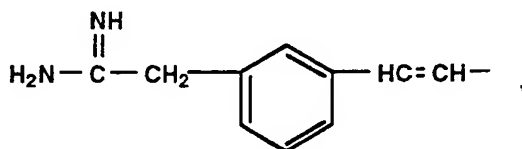
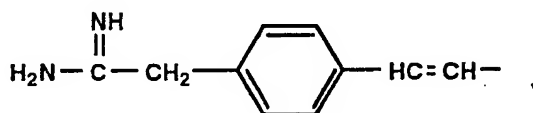
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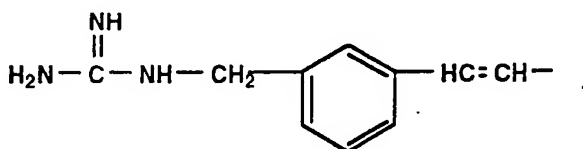
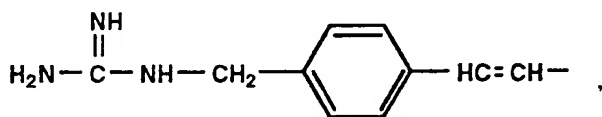
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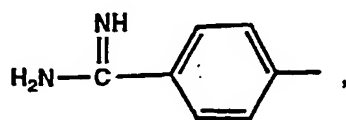
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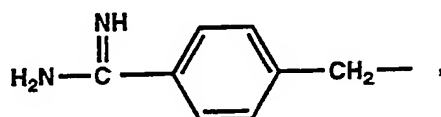
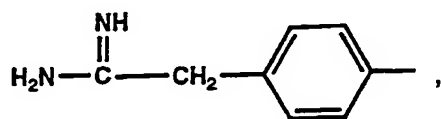
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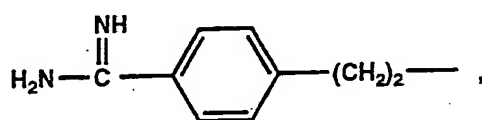
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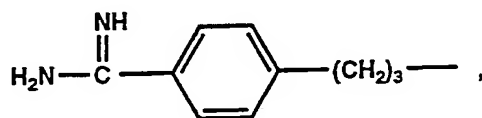
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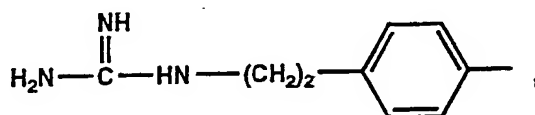
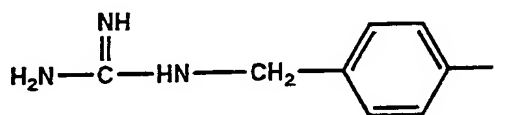
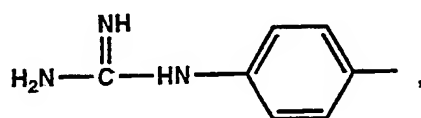
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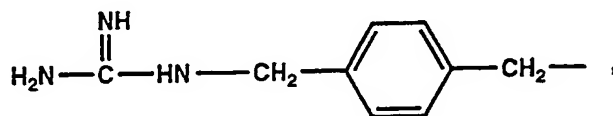
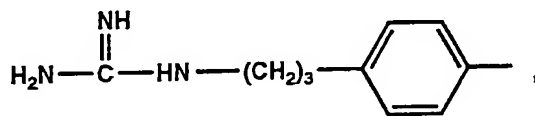
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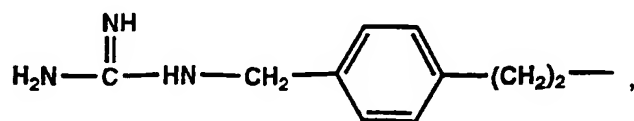


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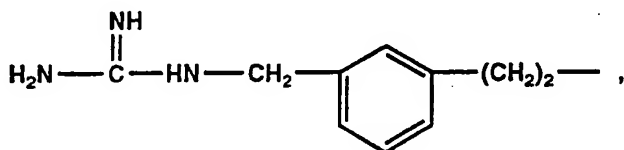
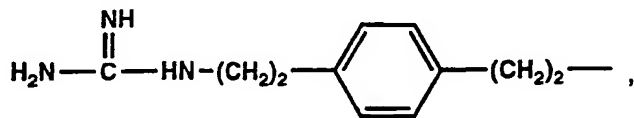


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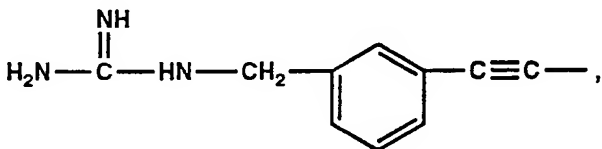
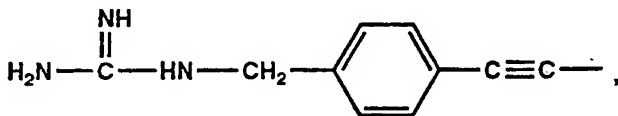




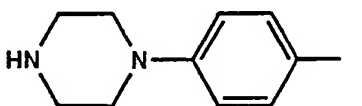
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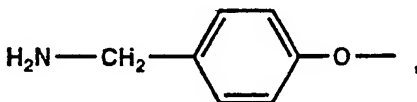
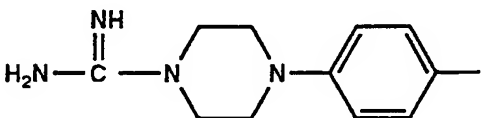
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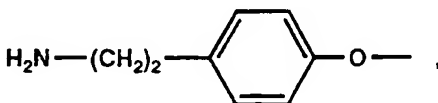
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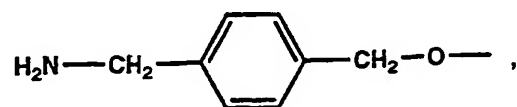
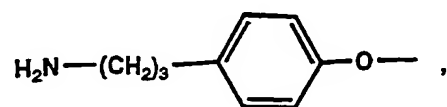


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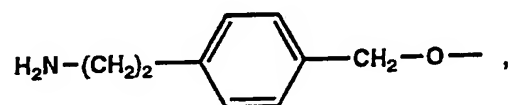
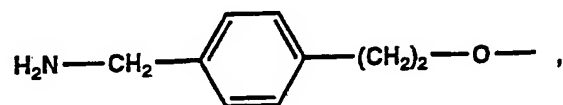


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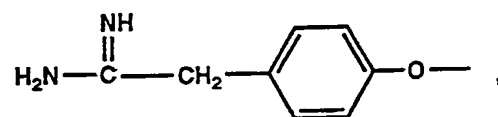




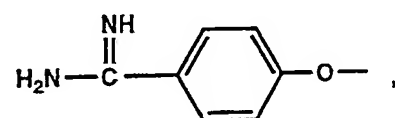
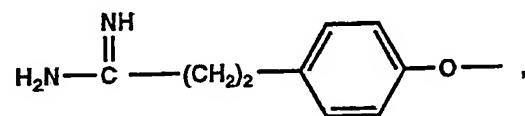
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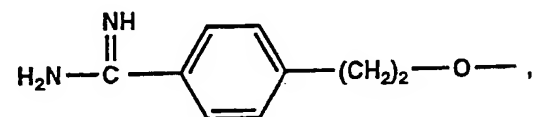
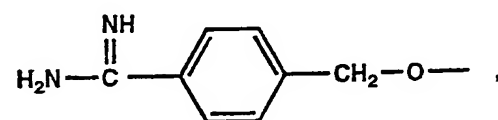
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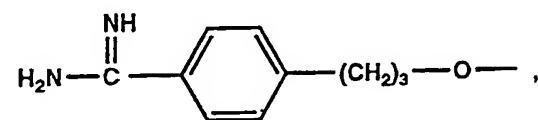
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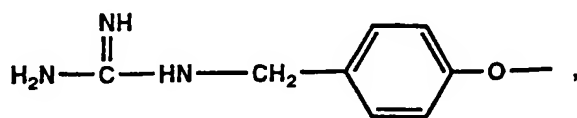
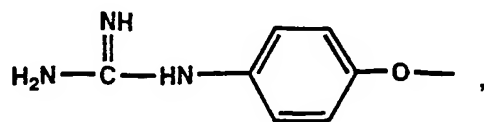
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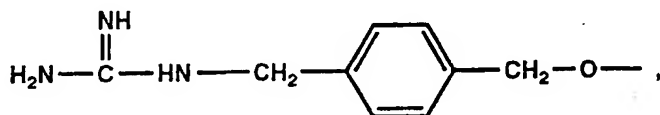
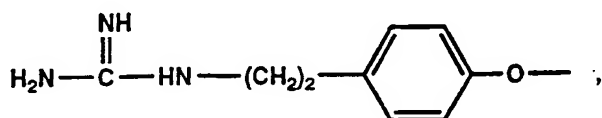
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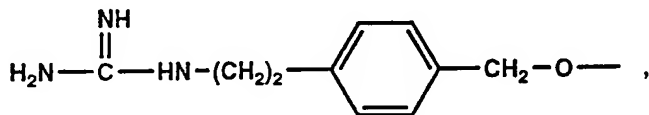
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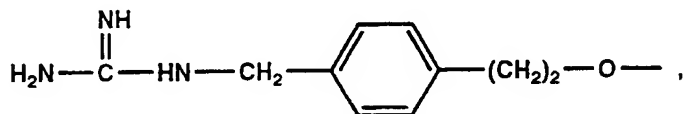
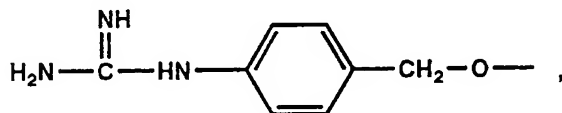
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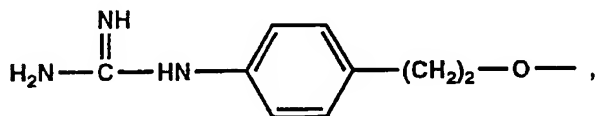
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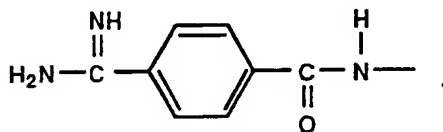
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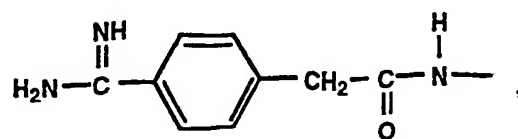


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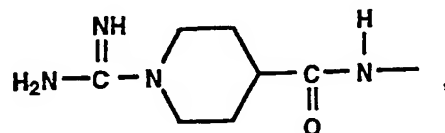


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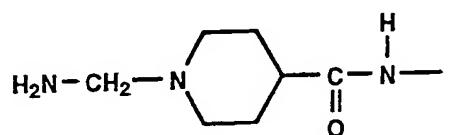




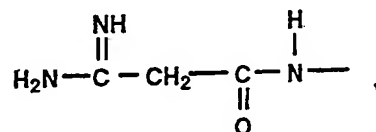
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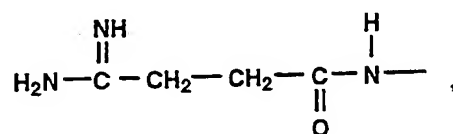
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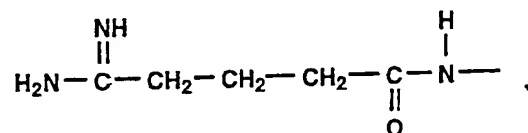
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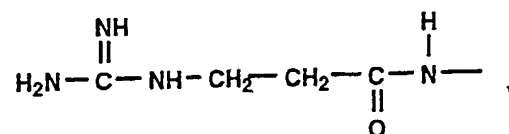
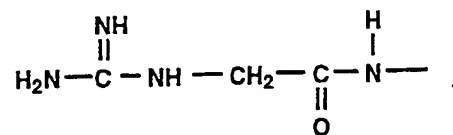
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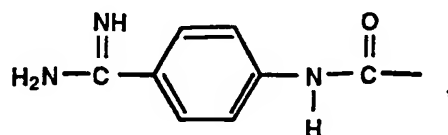


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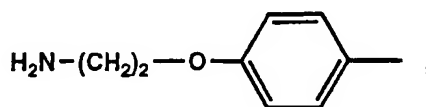


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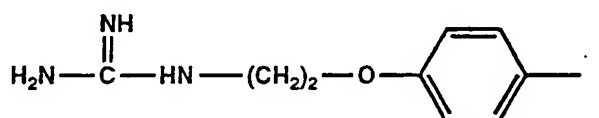
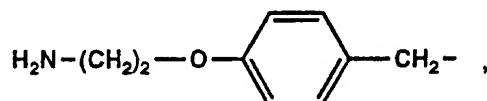




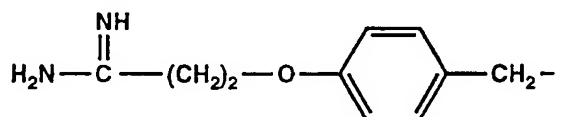
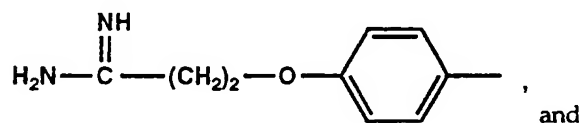
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6. The compound of Claim 2 selected from the group consisting of;

- 1-(4-chlorophenyl)methyl-4-(2-carboxyethyl)-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-  
 25 1,4-benzodiazapine-2,5-dione,  
 1-(diphenyl)methyl-4-(2-carboxyethyl)-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-  
 benzodiazapine-2,5-dione,  
 1-(*m*-trifluoromethyl)phenyl-4-(2-carboxyethyl)-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-  
 1H-1,4-benzodiazapine-2,5-dione,  
 30 1-methyl-4-(2-carboxyethyl)-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-  
 benzodiazapine-2,5-dione,  
 1-(1-methyl)ethyl-4-(2-carboxyethyl)-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-  
 benzodiazapine-2,5-dione,

- 1-(2-methyl)propyl-4-(2-carboxyethyl)-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(4-chlorophenyl)methyl-4-(2-carboxyethyl)-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
5 1-(diphenyl)methyl-4-(2-carboxyethyl)-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(*m*-trifluoromethyl)phenyl-4-(2-carboxyethyl)-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-methyl-4-(2-carboxyethyl)-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
10 1-(1-methyl)ethyl-4-(2-carboxyethyl)-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(2-methyl)propyl-4-(2-carboxyethyl)-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
15 1-(diphenyl)methyl-4-(2-carboxyethyl)-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(diphenyl)methyl-4-(2-carboxyethyl)-7-(6-aminohex-1-ynyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(diphenyl)methyl-4-(2-carboxyethyl)-7-(5-aminopentoxyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
20 1-(diphenyl)methyl-4-(2-carboxyethyl)-7-(5-guanidinopentyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(diphenyl)methyl-4-(2-carboxyethyl)-7-(4-guanidinobutoxy)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
25 1-(diphenyl)methyl-4-(2-carboxyethyl)-7-(5-aminopent-1-ynyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(diphenyl)methyl-4-(2-carboxyethyl)-7-(5-guanidinopent-1-ynyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(diphenyl)methyl-4-(2-carboxyethyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
30 1-(diphenyl)methyl-4-(2-carboxyethyl)-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(diphenyl)methyl-4-(2-carboxyethyl)-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
35 1-(diphenyl)methyl-4-(2-carboxyethyl)-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(4-chlorophenyl)methyl-4-(2-carboxyethyl)-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,



- 1-(4-chlorophenyl)methyl-4-(2-carboxyethyl)-7-(5-aminopentoxy)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(4-chlorophenyl)methyl-4-(2-carboxyethyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
5 1-(4-chlorophenyl)methyl-4-(2-carboxyethyl)-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(4-chlorophenyl)methyl-4-(2-carboxyethyl)-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(4-chlorophenyl)methyl-4-(2-carboxyethyl)-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
10 1-methyl-4-(2-carboxyethyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-methyl-4-(2-carboxyethyl)-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
15 1-methyl-4-(2-carboxyethyl)-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-methyl-4-(2-carboxyethyl)-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(1-methyl)ethyl-4-(2-carboxyethyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
20 1-(1-methyl)ethyl-4-(2-carboxyethyl)-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(1-methyl)ethyl-4-(2-carboxyethyl)-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
25 1-(2-methyl)ethyl-4-(2-carboxyethyl)-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(2-methyl)propyl-4-(2-carboxyethyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(2-methyl)propyl-4-(2-carboxyethyl)-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
30 1-(2-methyl)propyl-4-(2-carboxyethyl)-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(2-methyl)propyl-4-(2-carboxyethyl)-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
35 1-(2,4-difluorophenyl)methyl-4-(2-carboxyethyl)-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(2,4-difluorophenyl)methyl-4-(2-carboxyethyl)-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,

- 1-(2,4-difluorophenyl)methyl-4-(2-carboxyethyl)-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(2,4-difluorophenyl)methyl-4-(2-carboxyethyl)-7-(5-aminopentoxyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
5 1-(2,4-difluorophenyl)methyl-4-(2-carboxyethyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(2,4-difluorophenyl)methyl-4-(2-carboxyethyl)-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(2,4-difluorophenyl)methyl-4-(2-carboxyethyl)-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
10 1-(2,4-difluorophenyl)methyl-4-(2-carboxyethyl)-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(4-chlorophenyl)methyl-4-(2-carboxyethyl)-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
15 1-(diphenyl)methyl-4-(2-carboxyethyl)-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(*m*-trifluoromethyl)phenyl-4-(2-carboxyethyl)-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-methyl-4-(2-carboxyethyl)-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
20 1-(1-methyl)ethyl-4-(2-carboxyethyl)-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(2-methyl)propyl-4-(2-carboxyethyl)-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
25 1-(2,4-difluorophenyl)methyl-4-(2-carboxyethyl)-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(4-chlorophenyl)methyl-4-(3-butanoyl)-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(diphenyl)methyl-4-(3-butanoyl)-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
30 1-(*m*-trifluoromethyl)phenyl-4-(3-butanoyl)-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-methyl-4-(3-butanoyl)-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
35 1-(1-methyl)ethyl-4-(3-butanoyl)-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(2-methyl)propyl-4-(3-butanoyl)-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,

- 1-(4-chlorophenyl)methyl-4-(3-butanoyl)-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(diphenyl)methyl-4-(3-butanoyl)-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
5 1-(*m*-trifluoromethyl)phenyl-4-(3-butanoyl)-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-methyl-4-(3-butanoyl)-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(1-methyl)ethyl-4-(3-butanoyl)-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
10 1-(2-methyl)propyl-4-(3-butanoyl)-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(diphenyl)methyl-4-(3-butanoyl)-7-(6-aminoethyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
15 1-(diphenyl)methyl-4-(3-butanoyl)-7-(6-aminohex-1-ynyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(diphenyl)methyl-4-(3-butanoyl)-7-(5-aminopentoxyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(diphenyl)methyl-4-(3-butanoyl)-7-(5-guanidinopentyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
20 1-(diphenyl)methyl-4-(3-butanoyl)-7-(4-guanidinobutoxy)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(diphenyl)methyl-4-(3-butanoyl)-7-(5-aminopent-1-ynyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
25 1-(diphenyl)methyl-4-(3-butanoyl)-7-(5-guanidinopent-1-ynyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(diphenyl)methyl-4-(3-butanoyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(diphenyl)methyl-4-(3-butanoyl)-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
30 1-(diphenyl)methyl-4-(3-butanoyl)-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(diphenyl)methyl-4-(3-butanoyl)-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
35 1-(4-chlorophenyl)methyl-4-(3-butanoyl)-7-(6-aminoethyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(4-chlorophenyl)methyl-4-(3-butanoyl)-7-(5-aminopentoxyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,

- 1-(4-chlorophenyl)methyl-4-(3-butanoyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(4-chlorophenyl)methyl-4-(3-butanoyl)-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
5 1-(4-chlorophenyl)methyl-4-(3-butanoyl)-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(4-chlorophenyl)methyl-4-(3-butanoyl)-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-methyl-4-(3-butanoyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
10 dione,  
1-methyl-4-(3-butanoyl)-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-methyl-4-(3-butanoyl)-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
15 1-methyl-4-(3-butanoyl)-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(1-methyl)ethyl-4-(3-butanoyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(1-methyl)ethyl-4-(3-butanoyl)-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
20 1-(1-methyl)ethyl-4-(3-butanoyl)-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(2-methyl)ethyl-4-(3-butanoyl)-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
25 1-(2-methyl)propyl-4-(3-butanoyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(2-methyl)propyl-4-(3-butanoyl)-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(2-methyl)propyl-4-(3-butanoyl)-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
30 1-(2-methyl)propyl-4-(3-butanoyl)-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(2,4-difluorophenyl)methyl-4-(3-butanoyl)-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
35 1-(2,4-difluorophenyl)methyl-4-(3-butanoyl)-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(2,4-difluorophenyl)methyl-4-(3-butanoyl)-7-(6-aminoethyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,

- 1-(2,4-difluorophenyl)methyl-4-(3-butanoyl)-7-(5-aminopentoxy)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 1-(2,4-difluorophenyl)methyl-4-(3-butanoyl)-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 5 1-(2,4-difluorophenyl)methyl-4-(3-butanoyl)-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 1-(2,4-difluorophenyl)methyl-4-(3-butanoyl)-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 1-(2,4-difluorophenyl)methyl-4-(3-butanoyl)-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 10 1-(4-chlorophenyl)methyl-4-(3-butanoyl)-7-[2-(4-piperidinyloxy)]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 1-(diphenyl)methyl-4-(3-butanoyl)-7-[2-(4-piperidinyloxy)]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 15 1-(*m*-trifluoromethyl)phenyl-4-(3-butanoyl)-7-[2-(4-piperidinyloxy)]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 1-methyl-4-(3-butanoyl)-7-[2-(4-piperidinyloxy)]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 1-(1-methyl)ethyl-4-(3-butanoyl)-7-[2-(4-piperidinyloxy)]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 20 1-(2-methyl)propyl-4-(3-butanoyl)-7-[2-(4-piperidinyloxy)]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 1-(2,4-difluorophenyl)methyl-4-(3-butanoyl)-7-[2-(4-piperidinyloxy)]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 25 1-(4-chlorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 1-(diphenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 1-(*m*-trifluoromethyl)phenyl-4-[(2-methyl)carboxyethyl]-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 30 1-methyl-4-[(2-methyl)carboxyethyl]-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 1-(1-methyl)ethyl-4-[(2-methyl)carboxyethyl]-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 35 1-(2-methyl)propyl-4-[(2-methyl)carboxyethyl]-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 1-(4-chlorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,

- 1-(diphenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(*m*-trifluoromethyl)phenyl-4-[(2-methyl)carboxyethyl]-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
5 1-methyl-4-[(2-methyl)carboxyethyl]-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(1-methyl)ethyl-4-[(2-methyl)carboxyethyl]-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(2-methyl)propyl-4-[(2-methyl)carboxyethyl]-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
10 1-(diphenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(diphenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(6-aminohex-1-ynyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
15 1-(diphenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(5-aminopentoxyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(diphenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(5-guanidinopentyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(diphenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(4-guanidinobutoxy)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
20 1-(diphenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(5-aminopent-1-ynyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(diphenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(5-guanidinopent-1-ynyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
25 1-(diphenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(diphenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(diphenyl)methyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
30 1-(diphenyl)methyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(4-chlorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
35 1-(4-chlorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(5-aminopentoxyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(4-chlorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,

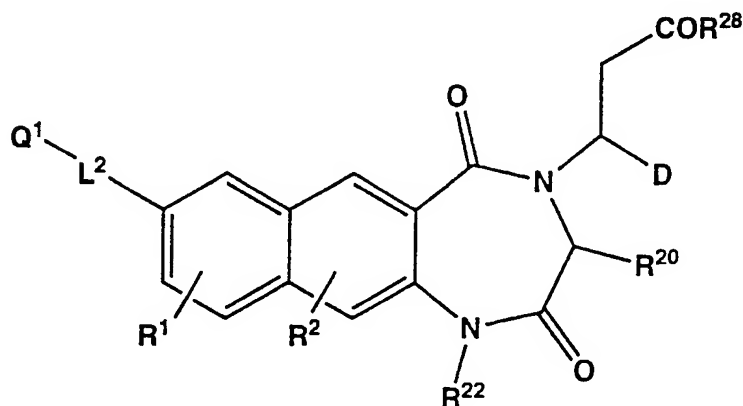
- 1-(4-chlorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 1-(4-chlorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 5 1-(4-chlorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 1-methyl-4-[(2-methyl)carboxyethyl]-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 1-methyl-4-[(2-methyl)carboxyethyl]-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 10 1-methyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 1-methyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 15 1-(1-methyl)ethyl-4-[(2-methyl)carboxyethyl]-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 1-(1-methyl)ethyl-4-[(2-methyl)carboxyethyl]-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 1-(1-methyl)ethyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 20 1-(2-methyl)ethyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 1-(2-methyl)propyl-4-[(2-methyl)carboxyethyl]-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 25 1-(2-methyl)propyl-4-[(2-methyl)carboxyethyl]-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 1-(2-methyl)propyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 1-(2-methyl)propyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 30 1-(2,4-difluorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(4-(1-piperizin)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 1-(2,4-difluorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(4-(2-aminoethoxy)-phenyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 35 1-(2,4-difluorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 1-(2,4-difluorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(5-aminopentoxyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,

- 1-(2,4-difluorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(4-amidino)benzamido-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(2,4-difluorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-(4-amidino)benzyloxy-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 5 1-(2,4-difluorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-amidinophenyl)]ethenyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(2,4-difluorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-amidinophenyl)]ethynyl-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 10 1-(4-chlorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(diphenyl)methyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(*m*-trifluoromethyl)phenyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 15 1-methyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(1-methyl)ethyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(2-methyl)propyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 20 1-(2,4-difluorophenyl)methyl-4-[(2-methyl)carboxyethyl]-7-[2-(4-piperidinyl)ethyloxy]-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(3-(butyryl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 25 1-(3-(butyryl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(5-guanidinopentyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(3-(4-aminobutyryl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(3-(4-guanidinobutyryl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(5-guanidinopentyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 30 1-(3-(6-aminohexoyl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(3-(6-guanidinoheptoyl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(5-guanidinopentyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,
- 35 1-(3-(6-aminopentoyl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(6-aminohexyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
1-(3-(6-guanidinopentoyl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(5-guanidinopentyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,

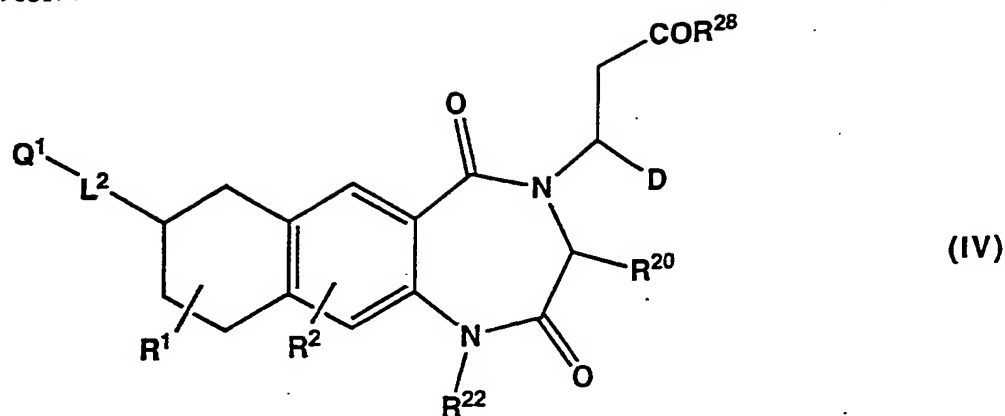


- 1-(3-(6-aminopropionyl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(6-aminoethyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
 1-(3-(6-guanidinopropionyl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(5-guanidinopentyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
 5 1-(3-(6-aminoacetyl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(6-aminoethyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
 1-(3-(6-guanidinoacetyl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(5-guanidinopentyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
 1-(3-(4-guanidinobenzoyl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(6-aminoethyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
 10 1-(4-(6-aminopropionyl)aminophenyl)methyl-4-(2-carboxyethyl)-7-(6-aminoethyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione,  
 1-(3-(4-aminobutyl)phenyl)methyl-4-(2-carboxyethyl)-7-(6-aminoethyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione, and  
 15 1-(4-carboxyphenyl)methyl-4-(2-carboxyethyl)-7-(6-aminoethyl)-3,4-dihydro-1H-1,4-benzodiazapine-2,5-dione.

- 20 7. The compound of Claim 1 represented by formula III and IV:



(III)



where

5  $R^1$  and  $R^2$  are independently selected from hydrogen and halo(F, Cl, Br, I);

$Q^1$  is selected from the group consisting of

- 10 (A) an amino group,  
 (B) an amidino group,  
 (C) an aminoalkyleneimino group,  
 (D) an iminoalkyleneamino group, and  
 (E) a guanidino group;

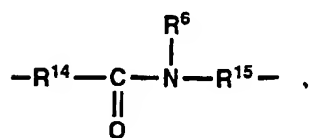
$R^{20}$  is hydrogen;

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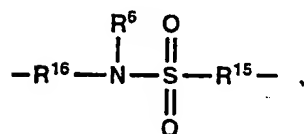
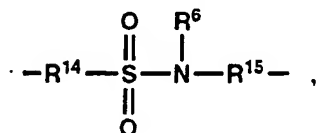
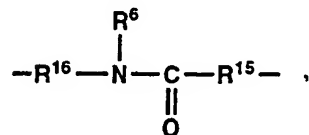
$R^{22}$  is selected from the group consisting of (i) hydrogen, (ii) optionally substituted C<sub>1</sub>-C<sub>12</sub>-alkyl, (iii) optionally substituted C<sub>6</sub>-C<sub>14</sub>-aryl, (iv) optionally substituted C<sub>3</sub>-C<sub>14</sub>-cycloalkyl, (v) optionally substituted C<sub>1</sub>-C<sub>12</sub>-alkyl-C<sub>6</sub>-C<sub>14</sub>-aryl, (vi) optionally substituted C<sub>1</sub>-C<sub>12</sub>-alkyl-C<sub>3</sub>-C<sub>14</sub>-cycloalkyl, where the substituents are selected from (a) halo (F, Cl, Br, I), (b) nitro, (c) hydroxy, (d) carboxy, (e) tetrazole, (f) hydroxamate, (g) sulfonamide, (h) trifluoroimide, (i) phosphonate, (j) C<sub>1</sub>-C<sub>6</sub>-alkyl, (k) C<sub>6</sub>-C<sub>14</sub>-aryl, (l) benzyl, (m) C<sub>3</sub>-C<sub>14</sub>-cycloalkyl, (n) COR<sup>24</sup> where R<sup>24</sup> is selected from the group C<sub>1</sub>-C<sub>8</sub>-alkoxy, C<sub>3</sub>-C<sub>12</sub>-alkenoxy, C<sub>6</sub>-C<sub>12</sub>-aryloxy, di-C<sub>1</sub>-C<sub>8</sub>-alkylamino-C<sub>1</sub>-C<sub>8</sub>-alkoxy, acylamino-C<sub>1</sub>-C<sub>8</sub>-alkoxy selected from the group acetylaminooethoxy, nicotinoylaminoethoxy, succinamidoethoxy, and pivaloyloxyethoxy, and C<sub>6</sub>-C<sub>12</sub>-aryl-C<sub>1</sub>-C<sub>8</sub>-alkoxy where the aryl group is unsubstituted or substituted with one to three of the groups nitro, halo (F, Cl, Br, I), C<sub>1</sub>-C<sub>4</sub>-alkoxy, amino, hydroxy, hydroxy-C<sub>2</sub>-C<sub>8</sub>-alkoxy, and dihydroxy-C<sub>3</sub>-C<sub>8</sub>-alkoxy, (o) CONR<sup>25</sup>R<sup>26</sup> where R<sup>25</sup> and R<sup>26</sup> are independently selected from hydrogen, C<sub>1</sub>-C<sub>10</sub>-alkyl, C<sub>3</sub>-C<sub>10</sub>-alkenyl, C<sub>6</sub>-C<sub>14</sub>-aryl, C<sub>1</sub>-C<sub>6</sub>-alkyl-C<sub>6</sub>-C<sub>10</sub>-aryl, optionally R<sup>25</sup> and R<sup>26</sup> taken together may form trimethylene, tetramethylene, pentamethylene, and 3-oxopentamethylene, and (vii) Q<sup>2</sup>-L<sup>3</sup> where Q<sup>2</sup> is selected from hydrogen, and Q<sup>1</sup>, and L<sup>3</sup> is selected from a chemical bond, and L<sup>1</sup>, where L<sup>1</sup> is an optionally substituted bivalent radical selected from the group consisting of C<sub>3</sub>-C<sub>7</sub>-alkylene, C<sub>3</sub>-C<sub>7</sub>-cycloalkylene, C<sub>3</sub>-C<sub>7</sub>-alkenylene, C<sub>4</sub>-C<sub>7</sub>-cycloalkenylene, C<sub>5</sub>-C<sub>8</sub>-cycloalkadienylene, C<sub>3</sub>-C<sub>7</sub>-alkadienylene, C<sub>3</sub>-C<sub>7</sub>-alkynylene, C<sub>4</sub>-C<sub>7</sub>-alkenynylene, C<sub>6</sub>-C<sub>14</sub>-arylene, C<sub>6</sub>-C<sub>14</sub>-aryl-C<sub>2</sub>-C<sub>4</sub>-alkynylene, C<sub>1</sub>-C<sub>3</sub>-alkyl-C<sub>6</sub>-C<sub>14</sub>-aryl-C<sub>2</sub>-C<sub>4</sub>-alkynylene, C<sub>6</sub>-C<sub>14</sub>-aryl-C<sub>2</sub>-C<sub>4</sub>-alkenylene, C<sub>1</sub>-C<sub>3</sub>-alkyl-C<sub>6</sub>-C<sub>14</sub>-arylene, C<sub>1</sub>-C<sub>3</sub>-alkyl-C<sub>6</sub>-C<sub>14</sub>-aryl-C<sub>2</sub>-C<sub>4</sub>-alkenylene, C<sub>6</sub>-C<sub>14</sub>-aryl-C<sub>1</sub>-C<sub>3</sub>-alkylene, C<sub>6</sub>-C<sub>14</sub>-aryl-C<sub>1</sub>-C<sub>3</sub>-alkyloxyene, C<sub>1</sub>-C<sub>2</sub>-alkyl-C<sub>6</sub>-C<sub>14</sub>-aryl-C<sub>1</sub>-C<sub>2</sub>-alkylene, C<sub>1</sub>-C<sub>3</sub>-alkyloxy-C<sub>6</sub>-C<sub>14</sub>-arylene, C<sub>2</sub>-C<sub>8</sub>-alkyloxyene, C<sub>1</sub>-C<sub>5</sub>-alkyloxy-

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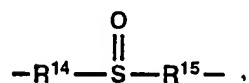
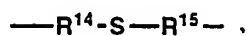
C<sub>1</sub>-C<sub>5</sub>-alkylene, C<sub>6</sub>-C<sub>10</sub>-aryloxyene, C<sub>6</sub>-C<sub>10</sub>-aryloxy-C<sub>1</sub>-C<sub>5</sub>-alkylene, C<sub>6</sub>-C<sub>10</sub>-arylthio-C<sub>1</sub>-C<sub>5</sub>-alkylene,



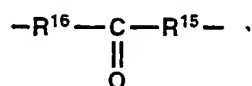
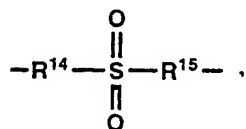
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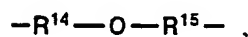
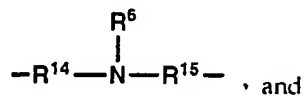
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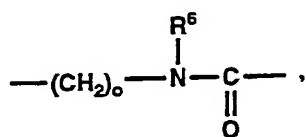
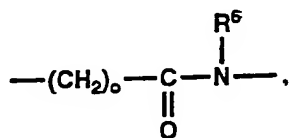
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where R<sup>14</sup> is selected from a chemical bond, C<sub>1</sub>-C<sub>8</sub>-alkyl, C<sub>3</sub>-C<sub>7</sub>-cycloalkyl, C<sub>2</sub>-C<sub>5</sub>-alkenyl, C<sub>3</sub>-C<sub>5</sub>-alkynyl, C<sub>6</sub>-C<sub>10</sub>-aryl, C<sub>1</sub>-C<sub>3</sub>-alkyl-C<sub>6</sub>-C<sub>12</sub>-aryl, C<sub>1</sub>-C<sub>2</sub>-alkyl-C<sub>6</sub>-C<sub>10</sub>-aryl-C<sub>1</sub>-C<sub>2</sub>-alkyl, and C<sub>6</sub>-C<sub>10</sub>-aryl-C<sub>1</sub>-C<sub>2</sub>-alkyl, R<sup>15</sup> is selected from a chemical bond, C<sub>1</sub>-C<sub>4</sub>-alkyl, C<sub>2</sub>-C<sub>4</sub>-alkenyl, C<sub>2</sub>-C<sub>4</sub>-alkynyl, and C<sub>6</sub>-C<sub>10</sub>-aryl, R<sup>16</sup> is selected from a chemical bond, C<sub>2</sub>-C<sub>5</sub>-alkyl, C<sub>3</sub>-C<sub>7</sub>-cycloalkyl, C<sub>3</sub>-C<sub>5</sub>-alkenyl, C<sub>3</sub>-C<sub>5</sub>-alkynyl, C<sub>6</sub>-C<sub>10</sub>-aryl, C<sub>1</sub>-C<sub>3</sub>-alkyl-C<sub>6</sub>-C<sub>12</sub>-aryl, C<sub>1</sub>-C<sub>3</sub>-alkyl-C<sub>6</sub>-C<sub>12</sub>-aryl,

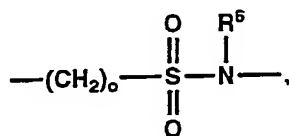
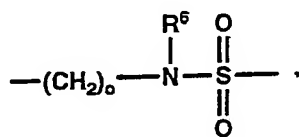
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and C<sub>6</sub>-C<sub>10</sub>-aryl-C<sub>1</sub>-C<sub>2</sub>-alkyl, where R<sup>14</sup> or R<sup>16</sup> is bonded to N and R<sup>15</sup> is bonded to Q<sup>2</sup> and where the substituents are selected from one to three R<sup>12</sup>;

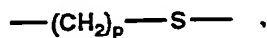
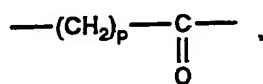
5 L<sup>2</sup> is an optionally substituted bivalent radical selected from the group consisting of



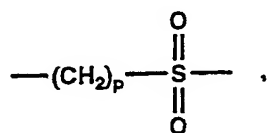
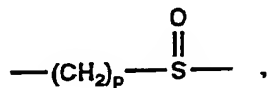
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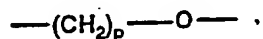
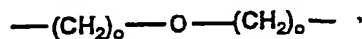
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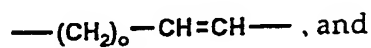
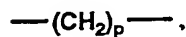
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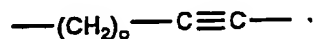
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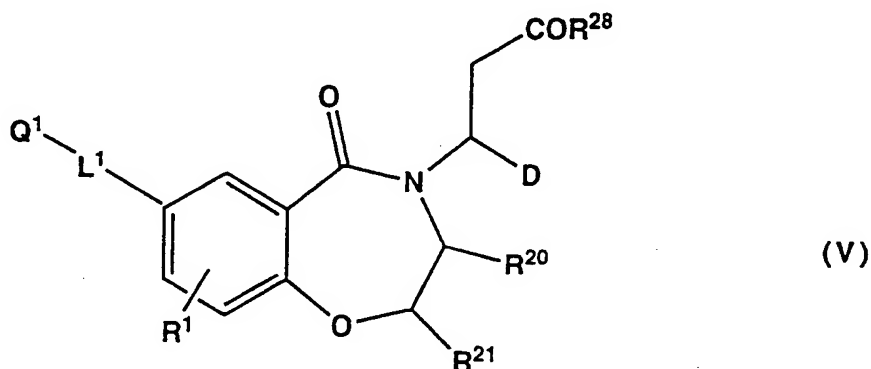
where o is 1 or 2, p is 1, 2, 3 or 4, R<sup>6</sup> is defined above, and where the substituents are selected from one to three R<sup>12</sup>;

- 5 D is selected from the group consisting of R<sup>22</sup>, and -(C=O)-Xaa, where Xaa is one to three D or L  $\alpha$ -amino acid;

- 10 R<sup>28</sup> is selected from the group consisting of (i) hydroxy, (ii) C<sub>1</sub>-C<sub>8</sub>-alkoxy, (iii) C<sub>3</sub>-C<sub>12</sub>-alkenoxy, (iv) C<sub>6</sub>-C<sub>12</sub>-aryloxy, (v) C<sub>1</sub>-C<sub>6</sub>-alkyl-C<sub>6</sub>-C<sub>12</sub>-aryloxy, (vi) di-C<sub>1</sub>-C<sub>8</sub>-alkylamino-C<sub>1</sub>-C<sub>8</sub>-alkoxy, (vii) acylamino-C<sub>1</sub>-C<sub>8</sub>-alkoxy selected from the group (a) acetylaminoethoxy, (b) nicotinoylaminoethoxy, and (c) succinamidoethoxy, (viii) C<sub>1</sub>-C<sub>8</sub>-alkoxyloxy-C<sub>1</sub>-C<sub>8</sub>-alkoxy, (ix) C<sub>6</sub>-C<sub>12</sub>-aryl-C<sub>1</sub>-C<sub>8</sub>-alkoxy where the aryl group is unsubstituted or substituted with one to three of the groups (a) nitro, (b) halo (F, Cl, Br, I), (c) C<sub>1</sub>-C<sub>4</sub>-alkoxy, and (d) amino, (x) hydroxy-C<sub>2</sub>-C<sub>8</sub>-alkoxy, (xi) dihydroxy-C<sub>3</sub>-C<sub>8</sub>-alkoxy, and (xii) NR<sup>29</sup>R<sup>30</sup> where R<sup>29</sup> and R<sup>30</sup> are independently selected from the group (a) hydrogen, (b) C<sub>1</sub>-C<sub>8</sub>-alkyl, (c) C<sub>3</sub>-C<sub>8</sub>-alkenyl, (d) C<sub>6</sub>-C<sub>12</sub>-aryl where the aryl group is unsubstituted or substituted with one to three of the groups nitro, halo (F, Cl, Br, I), C<sub>1</sub>-C<sub>4</sub>-alkoxy, and amino, and (e) C<sub>6</sub>-C<sub>12</sub>-aryl-C<sub>1</sub>-C<sub>8</sub>-alkyl where the aryl group is unsubstituted or substituted with one to three of the groups nitro, halo (F, Cl, Br, I), and C<sub>1</sub>-C<sub>4</sub>-alkoxy; and
- 20

pharmaceutically acceptable salts thereof.

- 25 8. The compound of Claim 1 represented by formula V:



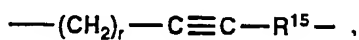
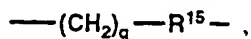
where

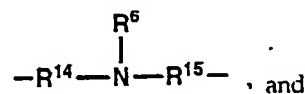
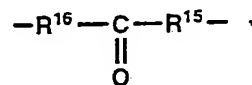
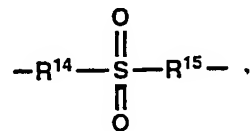
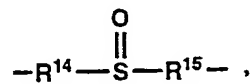
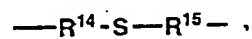
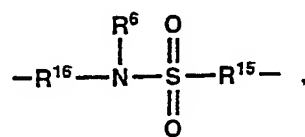
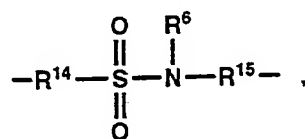
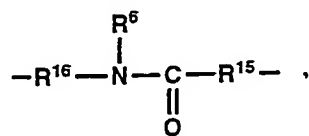
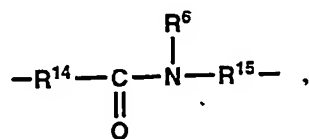
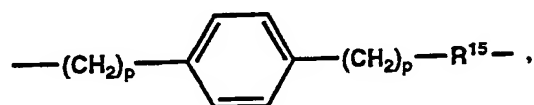
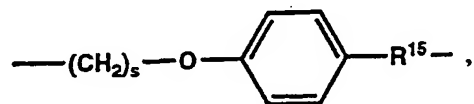
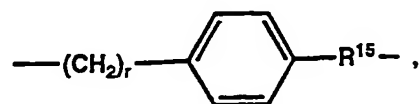
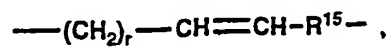
- 30 R<sup>1</sup> is one to three groups independently selected from hydrogen and halogen(F, Cl, Br, I);

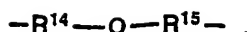
Q<sup>1</sup> is selected from the group consisting of

- 35 (A) an amino group,  
(B) an amidino group,  
(C) an aminoalkyleneimino group,  
(D) an iminoalkyleneamino group, and  
(E) a guanidino group;

- 40 L<sup>1</sup> is an optionally substituted bivalent radical selected from the group consisting of







where

5 p is 1, 2, 3, or 4;

q is 3, 4, 5, 6 or 7;

10 r is 1, 2, 3, 4, or 5;

s is 2 or 3;

R<sup>14</sup> is selected from

15 C<sub>2</sub>-C<sub>5</sub>-alkyl,  
C<sub>3</sub>-C<sub>7</sub>-cycloalkyl,  
C<sub>2</sub>-C<sub>5</sub>-alkenyl,  
C<sub>3</sub>-C<sub>5</sub>-alkynyl,  
C<sub>6</sub>-C<sub>10</sub>-aryl,  
C<sub>1</sub>-C<sub>2</sub>-alkyl-C<sub>6</sub>-C<sub>12</sub>-aryl,  
20 C<sub>1</sub>-C<sub>2</sub>-alkyl-C<sub>6</sub>-C<sub>10</sub>-aryl-C<sub>1</sub>-C<sub>2</sub>-alkyl,  
C<sub>6</sub>-C<sub>10</sub>-aryl-C<sub>1</sub>-C<sub>2</sub>-alkyl, and  
C<sub>6</sub>-C<sub>10</sub>-aryloxy-C<sub>1</sub>-C<sub>2</sub>-alkyl;

R<sup>15</sup> is a chemical bond connecting L<sup>1</sup> to position 8 of the benzazepine-one;

25

R<sup>16</sup> is selected from

30 C<sub>2</sub>-C<sub>5</sub>-alkyl,  
C<sub>3</sub>-C<sub>7</sub>-cycloalkyl,  
C<sub>3</sub>-C<sub>5</sub>-alkenyl,  
C<sub>3</sub>-C<sub>5</sub>-alkynyl,  
C<sub>6</sub>-C<sub>10</sub>-aryl,  
C<sub>1</sub>-C<sub>3</sub>-alkyl-C<sub>6</sub>-C<sub>12</sub>-aryl, and  
C<sub>6</sub>-C<sub>10</sub>-aryl-C<sub>1</sub>-C<sub>2</sub>-alkyl,

wherein R<sup>14</sup> and R<sup>16</sup> bond L<sup>1</sup> to Q<sup>1</sup>, and

35 where the substituents are selected from the group hydrogen, halo(F, Cl, Br, I), C<sub>1</sub>-C<sub>4</sub> alkoxy, C<sub>1</sub>-C<sub>4</sub> alkyl, phenyl, benzyl, and halo(F, Cl, Br, I)-C<sub>1</sub>-C<sub>4</sub> alkyl;

R<sup>20</sup> is hydrogen;

40 R<sup>21</sup> is selected from the group consisting of (i) hydrogen, (ii) optionally substituted C<sub>1</sub>-C<sub>12</sub>-alkyl, (iii) optionally substituted C<sub>6</sub>-C<sub>14</sub>-aryl, (iv) optionally substituted C<sub>3</sub>-C<sub>14</sub>-cycloalkyl, (v) optionally substituted C<sub>1</sub>-C<sub>12</sub>-alkyl-C<sub>6</sub>-C<sub>14</sub>-aryl, (vi) optionally substituted C<sub>1</sub>-C<sub>12</sub>-alkyl-C<sub>3</sub>-C<sub>14</sub>-cycloalkyl, where the substituents are selected from (a) halo (F, Cl, Br, I), (b) nitro, (c) hydroxy, (d) carboxy, (e) tetrazole, (f) hydroxamate, (g) sulfonamide, (h) trifluoroimide, (i) phosphonate, (j) C<sub>1</sub>-C<sub>6</sub>-alkyl, (k) C<sub>6</sub>-C<sub>14</sub>-aryl, (l) benzyl, (m) C<sub>3</sub>-C<sub>14</sub>-cycloalkyl, (n) COR<sup>24</sup> where R<sup>24</sup> is selected from the group C<sub>1</sub>-C<sub>8</sub>-alkoxy, C<sub>3</sub>-C<sub>12</sub>-alkenoxy, C<sub>6</sub>-C<sub>12</sub>-aryloxy, di-C<sub>1</sub>-C<sub>8</sub>-alkylamino-C<sub>1</sub>-C<sub>8</sub>-alkoxy, acylamino-C<sub>1</sub>-C<sub>8</sub>-alkoxy selected from the group acetylaminoethoxy, nicotinoylaminoethoxy, succinamidoethoxy, and pivaloyloxyethoxy, and C<sub>6</sub>-C<sub>12</sub>-aryl-C<sub>1</sub>-C<sub>8</sub>-alkoxy where the aryl group is unsubstituted or substituted with one to three of the groups nitro, halo (F, Cl, Br, I), C<sub>1</sub>-C<sub>4</sub>-alkoxy, amino, hydroxy, hydroxy-C<sub>2</sub>-C<sub>8</sub>-alkoxy, and dihydroxy-C<sub>3</sub>-C<sub>8</sub>-alkoxy, (o) CONR<sup>25</sup>R<sup>26</sup> where R<sup>25</sup> and R<sup>26</sup> are independently selected from hydrogen, C<sub>1</sub>-C<sub>10</sub>-alkyl, C<sub>3</sub>-C<sub>10</sub>-alkenyl, C<sub>6</sub>-C<sub>14</sub>-aryl, C<sub>1</sub>-C<sub>6</sub>-alkyl-C<sub>6</sub>-C<sub>10</sub>-aryl,

50

optionally  $R^{25}$  and  $R^{26}$  taken together may form trimethylene, tetramethylene, pentamethylene, and 3-oxopentamethylene, and (vii)  $Q^2-L^3$  where  $Q^2$  is selected from hydrogen and  $Q^1$ , and  $L^3$  is selected from a chemical bond and  $L^1$  where  $R^{15}$  bonds  $L^1$  to position 4 of the benzazepine-one;

5

D is selected from the group consisting of  $R^{21}$ , and  $-(C=O)-Xaa$ , where Xaa is one to three D or L  $\alpha$ -amino acid;

10

$R^{28}$  is selected from the group consisting of (i) hydroxy, (ii)  $C_1$ -C<sub>8</sub>-alkoxy, (iii)  $C_3$ -C<sub>12</sub>-alkenoxy, (iv)  $C_6$ -C<sub>12</sub>-aryloxy, (v)  $C_1$ -C<sub>6</sub>-alkyl- $C_6$ -C<sub>12</sub>-aryloxy, (vi) di- $C_1$ -C<sub>8</sub>-alkylamino- $C_1$ -C<sub>8</sub>-alkoxy, (vii) acylamino- $C_1$ -C<sub>8</sub>-alkoxy selected from the group (a) acetylaminoethoxy, (b) nicotinoylaminoethoxy, and (c) succinamidoethoxy, (viii)  $C_1$ -C<sub>8</sub>-alkoxyloxy- $C_1$ -C<sub>8</sub>-alkoxy, (ix)  $C_6$ -C<sub>12</sub>-aryl- $C_1$ -C<sub>8</sub>-alkoxy where the aryl group is unsubstituted or substituted with one to three of the groups (a) nitro, (b) halo (F, Cl, Br, I), (c)  $C_1$ -C<sub>4</sub>-alkoxy, and (d) amino, (x) hydroxy- $C_2$ -C<sub>8</sub>-alkoxy, (xi) dihydroxy- $C_3$ -C<sub>8</sub>-alkoxy, and (xii)  $NR^{29}R^{30}$  where  $R^{29}$  and  $R^{30}$  are independently selected from the group (a) hydrogen, (b)  $C_1$ -C<sub>8</sub>-alkyl, (c)  $C_3$ -C<sub>8</sub>-alkenyl, (d)  $C_6$ -C<sub>12</sub>-aryl where the aryl group is unsubstituted or substituted with one to three of the groups nitro, halo (F, Cl, Br, I),  $C_1$ -C<sub>4</sub>-alkoxy, and amino, and (e)  $C_6$ -C<sub>12</sub>-aryl- $C_1$ -C<sub>8</sub>-alkyl where the aryl group is unsubstituted or substituted with one to three of the groups nitro, halo (F, Cl, Br, I), and  $C_1$ -C<sub>4</sub>-alkoxy; and

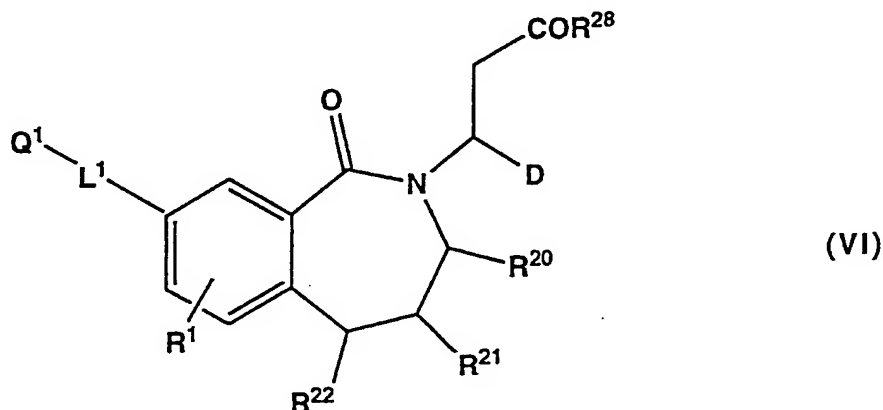
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20

pharmaceutically acceptable salts thereof.

25

9. The compound of Claim 1 represented by formula VI:



30

where

$R^1$  is one to three groups independently selected from hydrogen and halogen(F, Cl, Br, I);

35

$Q^1$  is selected from the group consisting of

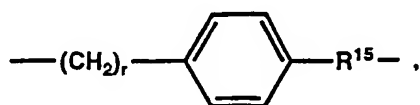
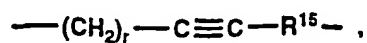
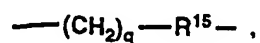
- (A) an amino group,
- (B) an amidino group,
- (C) an aminoalkyleneimino group,
- (D) an iminoalkyleneamino group, and
- (E) a guanidino group;

40

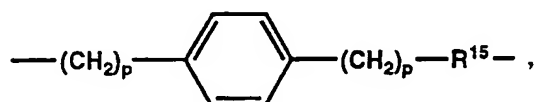
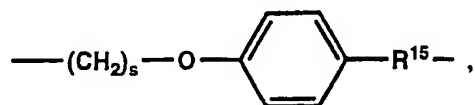
$L^1$  is an optionally substituted bivalent radical selected from the group consisting of



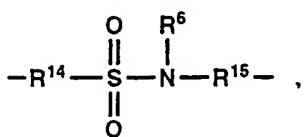
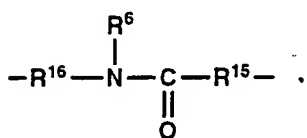
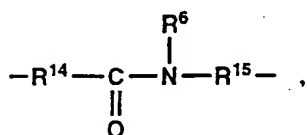
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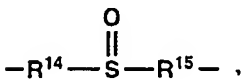
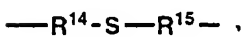
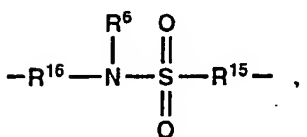
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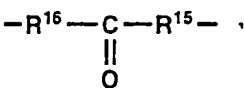
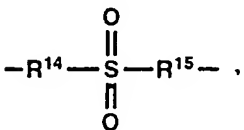
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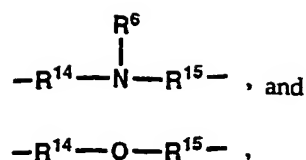


20



25





5 where

p is 1, 2, 3, or 4;

q is 3, 4, 5, 6 or 7;

10

r is 1, 2, 3, 4, or 5;

s is 2 or 3;

15

R<sup>14</sup> is selected from

C<sub>2</sub>-C<sub>5</sub>-alkyl,

C<sub>3</sub>-C<sub>7</sub>-cycloalkyl,

C<sub>2</sub>-C<sub>5</sub>-alkenyl,

C<sub>3</sub>-C<sub>5</sub>-alkynyl,

20

C<sub>6</sub>-C<sub>10</sub>-aryl,

C<sub>1</sub>-C<sub>2</sub>-alkyl-C<sub>6</sub>-C<sub>12</sub>-aryl,

C<sub>1</sub>-C<sub>2</sub>-alkyl-C<sub>6</sub>-C<sub>10</sub>-aryl-C<sub>1</sub>-C<sub>2</sub>-alkyl,

C<sub>6</sub>-C<sub>10</sub>-aryl-C<sub>1</sub>-C<sub>2</sub>-alkyl, and

C<sub>6</sub>-C<sub>10</sub>-aryloxy-C<sub>1</sub>-C<sub>2</sub>-alkyl;

25

R<sup>15</sup> is a chemical bond connecting L<sup>1</sup> to position 8 of the benzazepine-one;

R<sup>16</sup> is selected from

C<sub>2</sub>-C<sub>5</sub>-alkyl,

30

C<sub>3</sub>-C<sub>7</sub>-cycloalkyl,

C<sub>3</sub>-C<sub>5</sub>-alkenyl,

C<sub>3</sub>-C<sub>5</sub>-alkynyl,

C<sub>6</sub>-C<sub>10</sub>-aryl,

C<sub>1</sub>-C<sub>3</sub>-alkyl-C<sub>6</sub>-C<sub>12</sub>-aryl, and

35

C<sub>6</sub>-C<sub>10</sub>-aryl-C<sub>1</sub>-C<sub>2</sub>-alkyl,

wherein R<sup>14</sup> and R<sup>16</sup> bond L<sup>1</sup> to Q<sup>1</sup>, and

where the substituents are selected from the group hydrogen, halo(F, Cl, Br, I), C<sub>1</sub>-

C<sub>4</sub> alkoxy, C<sub>1</sub>-C<sub>4</sub> -alkyl, phenyl, benzyl, and halo(F, Cl, Br, I)-C<sub>1</sub>-C<sub>4</sub>-alkyl;

40

R<sup>20</sup> is hydrogen;

R<sup>21</sup> and R<sup>22</sup> are selected from the group consisting of (i) hydrogen, (ii) optionally

substituted C<sub>1</sub>-C<sub>12</sub>-alkyl, (iii) optionally substituted C<sub>6</sub>-C<sub>14</sub>-aryl, (iv) optionally

45

substituted C<sub>3</sub>-C<sub>14</sub>-cycloalkyl, (v) optionally substituted C<sub>1</sub>-C<sub>12</sub>-alkyl-C<sub>6</sub>-C<sub>14</sub>-

aryl, (vi) optionally substituted C<sub>1</sub>-C<sub>12</sub>-alkyl-C<sub>3</sub>-C<sub>14</sub>-cycloalkyl, where the

substituents are selected from (a) halo (F, Cl, Br, I), (b) nitro, (c) hydroxy, (d)

carboxy, (e) tetrazole, (f) hydroxamate, (g) sulfonamide, (h) trifluoroimide, (i)

phosphonate, (j) C<sub>1</sub>-C<sub>6</sub>-alkyl, (k) C<sub>6</sub>-C<sub>14</sub>-aryl, (l) benzyl, (m) C<sub>3</sub>-C<sub>14</sub>-cycloalkyl,

50

(n) COR<sup>24</sup> where R<sup>24</sup> is selected from the group C<sub>1</sub>-C<sub>8</sub>-alkoxy, C<sub>3</sub>-C<sub>12</sub>-alkenoxy,

C<sub>6</sub>-C<sub>12</sub>-aryloxy, di-C<sub>1</sub>-C<sub>8</sub>-alkylamino-C<sub>1</sub>-C<sub>8</sub>-alkoxy, acylamino-C<sub>1</sub>-C<sub>8</sub>-alkoxy

selected from the group acetylaminooethoxy, nicotinoylaminooethoxy,

5 succinamidoethoxy, and pivaloyloxyethoxy, and C<sub>6</sub>-C<sub>12</sub>-aryl-C<sub>1</sub>-C<sub>8</sub>-alkoxy where the aryl group is unsubstituted or substituted with one to three of the groups nitro, halo (F, Cl, Br, I), C<sub>1</sub>-C<sub>4</sub>-alkoxy, amino, hydroxy, hydroxy-C<sub>2</sub>-C<sub>8</sub>-alkoxy, and dihydroxy-C<sub>3</sub>-C<sub>8</sub>-alkoxy, (n) CONR<sup>25</sup>R<sup>26</sup> where R<sup>25</sup> and R<sup>26</sup> are independently selected from hydrogen, C<sub>1</sub>-C<sub>10</sub>-alkyl, C<sub>3</sub>-C<sub>10</sub>-alkenyl, C<sub>6</sub>-C<sub>14</sub>-aryl, C<sub>1</sub>-C<sub>6</sub>-alkyl-C<sub>6</sub>-C<sub>10</sub>-aryl, optionally R<sup>25</sup> and R<sup>26</sup> taken together may form trimethylene, tetramethylene, pentamethylene, and 3-oxopentamethylene, and (vii) Q<sup>2</sup>-L<sup>3</sup> where Q<sup>2</sup> is selected from hydrogen and Q<sup>1</sup>, and L<sup>3</sup> is selected from a chemical bond and L<sup>1</sup> where R<sup>15</sup> bonds L<sup>1</sup> to position 4 or 5 of the benzazepine-one;

10 D is selected from the group consisting of R<sup>21</sup>, and -(C=O)-Xaa, where Xaa is one to three D or L α-amino acid;

15 R<sup>28</sup> is selected from the group consisting of (i) hydroxy, (ii) C<sub>1</sub>-C<sub>8</sub>-alkoxy, (iii) C<sub>3</sub>-C<sub>12</sub>-alkenoxy, (iv) C<sub>6</sub>-C<sub>12</sub>-aryloxy, (v) C<sub>1</sub>-C<sub>6</sub>-alkyl-C<sub>6</sub>-C<sub>12</sub>-aryloxy, (vi) di-C<sub>1</sub>-C<sub>8</sub>-alkylamino-C<sub>1</sub>-C<sub>8</sub>-alkoxy, (vii) acylamino-C<sub>1</sub>-C<sub>8</sub>-alkoxy selected from the group (a) acetylaminooethoxy, (b) nicotinoylaminoethoxy, and (c) succinamidoethoxy, (viii) C<sub>1</sub>-C<sub>8</sub>-alkoxyloxy-C<sub>1</sub>-C<sub>8</sub>-alkoxy, (ix) C<sub>6</sub>-C<sub>12</sub>-aryl-C<sub>1</sub>-C<sub>8</sub>-alkoxy where the aryl group is unsubstituted or substituted with one to three of the groups (a) nitro, (b) halo (F, Cl, Br, I), (c) C<sub>1</sub>-C<sub>4</sub>-alkoxy, and (d) amino, (x) hydroxy-C<sub>2</sub>-C<sub>8</sub>-alkoxy, (xi) dihydroxy-C<sub>3</sub>-C<sub>8</sub>-alkoxy, and (xii) NR<sup>29</sup>R<sup>30</sup> where R<sup>29</sup> and R<sup>30</sup> are independently selected from the group (a) hydrogen, (b) C<sub>1</sub>-C<sub>8</sub>-alkyl, (c) C<sub>3</sub>-C<sub>8</sub>-alkenyl, (d) C<sub>6</sub>-C<sub>12</sub>-aryl where the aryl group is unsubstituted or substituted with one to three of the groups nitro, halo (F, Cl, Br, I), C<sub>1</sub>-C<sub>4</sub>-alkoxy, and amino, and (e) C<sub>6</sub>-C<sub>12</sub>-aryl-C<sub>1</sub>-C<sub>8</sub>-alkyl where the aryl group is unsubstituted or substituted with one to three of the groups nitro, halo (F, Cl, Br, I), and C<sub>1</sub>-C<sub>4</sub>-alkoxy; and

pharmaceutically acceptable salts thereof.

30 10. A pharmaceutical composition comprising a pharmaceutically acceptable excipient and the compound of Claims 1-9.

35 11. A method of inhibiting platelet aggregation comprising administering a platelet aggregation inhibiting amount of the composition of Claim 10.

12. A method for reducing platelet aggregation in a mammal comprising administering a pharmaceutically effective amount of the composition of Claim 10 to the mammal.

40 13. A method for treating a mammal having an increased propensity for thrombus formation comprising administering a pharmaceutically effective amount of the composition of Claim 10 to the mammal.

45 14. The method of Claim 13 further comprising administering the composition in combination with a thrombolytic agent.

15. The method of Claim 13 further comprising administering the composition in combination with an anticoagulant.

## INTERNATIONAL SEARCH REPORT

International Application No

PCT/US 92/08788

**I. CLASSIFICATION OF SUBJECT MATTER** (If several classification symbols apply, indicate all)<sup>6</sup>

According to International Patent Classification (IPC) or to both National Classification and IPC

Int.Cl. 5 C07D243/14; A61K31/55; C07D223/16; C07D267/14

**II. FIELDS SEARCHED**Minimum Documentation Searched<sup>7</sup>

Classification System	Classification Symbols
Int.Cl. 5	C07D

Documentation Searched other than Minimum Documentation  
to the Extent that such Documents are Included in the Fields Searched<sup>8</sup>**III. DOCUMENTS CONSIDERED TO BE RELEVANT<sup>9</sup>**

Category <sup>9</sup>	Citation of Document, <sup>11</sup> with indication, where appropriate, of the relevant passages <sup>12</sup>	Relevant to Claim No. <sup>13</sup>
A	WO,A,9 109 024 (KYORIN PHARMACEUTICAL CO.) 27 June 1991 & EP,A,0 456 835 cited in the application ---	1, 10
A	EP,A,0 268 148 (F. HOFFMANN-LA ROCHE AG) 25 May 1988 see example 40 -----	

<sup>9</sup> Special categories of cited documents: <sup>10</sup><sup>"A"</sup> document defining the general state of the art which is not considered to be of particular relevance<sup>"E"</sup> earlier document but published on or after the international filing date<sup>"L"</sup> document which may throw doubts on priority claim(s) or which is cited to establish the publication date of another citation or other special reason (as specified)<sup>"O"</sup> document referring to an oral disclosure, use, exhibition or other means<sup>"P"</sup> document published prior to the international filing date but later than the priority date claimed<sup>"T"</sup> later document published after the international filing date or priority date and not in conflict with the application but cited to understand the principle or theory underlying the invention<sup>"X"</sup> document of particular relevance; the claimed invention cannot be considered novel or cannot be considered to involve an inventive step<sup>"Y"</sup> document of particular relevance; the claimed invention cannot be considered to involve an inventive step when the document is combined with one or more other such documents, such combination being obvious to a person skilled in the art.<sup>"&"</sup> document member of the same patent family**IV. CERTIFICATION**

Date of the Actual Completion of the International Search

29 DECEMBER 1992

Date of Mailing of this International Search Report

28.01.93

International Searching Authority

EUROPEAN PATENT OFFICE

Signature of Authorized Officer

DE JONG B.S.

## INTERNATIONAL SEARCH REPORT

International application No.

PCT/US 92/08788

**Box I Observations where certain claims were found unsearchable (Continuation of item 1 of first sheet)**

This international search report has not been established in respect of certain claims under Article 17(2)(a) for the following reasons:

1. ☐ Claims Nos.:  
because they relate to subject matter not required to be searched by this Authority, namely:  
Remark: Although claims 11-15 are directed to a method of treatment of (diagnostic method practised on) the human/animal body the search has been carried out and based on the alleged effects of the compound/composition.
2. ☐ Claims Nos.:  
because they relate to parts of the international application that do not comply with the prescribed requirements to such an extent that no meaningful international search can be carried out, specifically:  
Claims not searched: 1  
Claim 1 is so broad that a meaningful search was not possible for economical reasons (Guidelines BIII, 2-1).
3. ☐ Claims Nos.:  
because they are dependent claims and are not drafted in accordance with the second and third sentences of Rule 6.4(a).

**Box II Observations where unity of invention is lacking (Continuation of item 2 of first sheet)**

This International Searching Authority found multiple inventions in this international application, as follows:

1. ☐ As all required additional search fees were timely paid by the applicant, this international search report covers all searchable claims.
2. ☐ As all searchable claims could be searched without effort justifying an additional fee, this Authority did not invite payment of any additional fee.
3. ☐ As only some of the required additional search fees were timely paid by the applicant, this international search report covers only those claims for which fees were paid, specifically claims Nos.:
4. ☐ No required additional search fees were timely paid by the applicant. Consequently, this international search report is restricted to the invention first mentioned in the claims; it is covered by claims Nos.:

Remark on Protest

☐ The additional search fees were accompanied by the applicant's protest.☐ No protest accompanied the payment of additional search fees.

**ANNEX TO THE INTERNATIONAL SEARCH REPORT  
ON INTERNATIONAL PATENT APPLICATION NO. US 9208788  
SA 66010**

This annex lists the patent family members relating to the patent documents cited in the above-mentioned international search report.  
The members are as contained in the European Patent Office EDP file on  
The European Patent Office is in no way liable for these particulars which are merely given for the purpose of information. 29/12/92

Patent document cited in search report	Publication date	Patent family member(s)	Publication date
WO-A-9109024	27-06-91	JP-A- 3181469	07-08-91
		AU-A- 6890591	18-07-91
		EP-A- 0456835	21-11-91
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EP-A-0268148	25-05-88	AU-B- 600769	23-08-90
		AU-A- 8090987	19-05-88
		DE-A- 3775177	23-01-92
		JP-A- 63139171	10-06-88
		US-A- 4808605	28-02-89
		ZA-A- 8708362	16-05-88
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